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INTERNATIONAL APPLICATION PUBLISHED UNDER THE PATENT COOPERATION TREATY (PCT)

(54) Title: MALTOGENIC ALPHA–AMYLASE VARIANTS

(57) Abstract

The inventors have modified the amino acid sequence of a maltogenic alpha–amylase to obtain variants with improved properties, based on the three–dimensional structure of the maltogenic alpha–amylase Novamyl. The variants have altered physicochemical properties., e.g. an altered pH optimum, improved thermostability, increased specific activity, an altered cleavage pattern or an increased ability to reduce retrogradation of starch or staling of bread.

MALTOGENIC ALPHA-AMYLASE VARIANTS

FIELD OF THE INVENTION

The present invention relates to variants of maltogenic amylase and to methods of constructing such variants.

5 **BACKGROUND OF THE INVENTION**

Maltogenic alpha-amylase (glucan 1,4-α-maltohydrolase, E.C. 3.2.1.133) is able to hydrolyze amylose and amylopectin to maltose in the alpha-configuration, and is also able to hydrolyze maltotriose as well as cyclodextrin.

A maltogenic alpha-amylase from *Bacillus* (EP 120 693) is commercially
10 available under the trade name Novamyl® (product of Novo Nordisk A/S, Denmark) and is widely used in the baking industry as an anti-staling agent due to its ability to reduce retrogradation of starch. Novamyl® shares several characteristics with cyclodextrin glucanotransferases (CGTases), including sequence homology (Henrissat B., Bairoch A. 1996) and formation of transglycosylation products (Christophersen, C., et al., 1997,
15 Starch, vol. 50, No. 1, 39-45).

Cyclomaltodextrin glucanotransferase (E.C. 2.4.1.19), also designated cyclodextrin glucanotransferase or cyclodextrin glycosyltransferase, abbreviated herein as CGTase, catalyses the conversion of starch and similar substrates into cyclomaltodextrins via an intramolecular transglycosylation reaction, thereby forming
20 cyclomaltodextrins (or CD) of various sizes.

CGTases are widely distributed and from several different bacterial sources, including *Bacillus*, *Brevibacterium*, *Clostridium*, *Corynebacterium*, *Klebsiella*, *Micrococcus*, *Thermoanaerobacter* and *Thermoanaerobacterium* have been extensively described in the literature. A CGTase produced by *Thermoanaerobacter sp.*
25 has been reported in *Norman B E, Jørgensen S T*; Denpun Kagaku 1992 39 99-106, and WO 89/03421, and the amino acid sequence has been disclosed in WO 96/33267. The sequence of CGTases from *Thermoanaerobacterium thermosulfurigenes* and from *Bacillus circulans* is available on the Internet (SCOP or PDF home pages) as pdf file 1CIU, and the sequence of a CGTase from *B. circulans* is available as pdf file 1CDG.
30 Tachibana, Y., Journal of Fermentation and Bioengineering, 83 (6), 540-548 (1997) describes the cloning and expression of a CGTase. Variants of CGTases have been described by Kim, Y. H., Biochemistry and Molecular Biology International, 41 (2), 227-234 (1997); Sin K-A, Journal of Biotechnology, 32 (3), 283-288 (1994); D Penninga, Biochemistry, 34 (10), 3368-3376 (1995); and WO 96/33267.

Recently, the tertiary structure of several CGTases have been reported. *Hofman et al.*
[*Hofman B E, Bender H, Schultz G E; J. Mol. Biol.* 1989 **209** 793-800] and *Klein &*
Schulz [*Klein C, Schulz G E; J. Mol. Biol.* 1991 **217** 737-750] report the tertiary
structure of a CGTase derived from *Bacillus circulans* Strain 8, *Kubota et al.* [*Kubota M,*
5 *Matsuura Y, Sakai S and Katsube Y; Denpun Kagaku* 1991 **38** 141-146] report the
tertiary structure of a CGTase derived from *Bacillus stearothermophilus* TC-91, *Lawson*
et al. [Lawson C L, van Montfort R, Strokopytov B, Rozeboom H J, Kalk K H, de Vries
G E, Penninga D, Dijkhuizen L, and Dijkstra B W; J. Mol. Biol. 1994 **236** 590-600]
report the tertiary structure of a CGTase derived from *Bacillus circulans* Strain 251,
10 *Strokopytov et al.* [*Strokopytov B, Penninga D, Rozeboom H J; Kalk K H, Dijkhuizen L*
and Dijkstra B W; Biochemistry 1995 **34** 2234-2240] report the tertiary structure of a
CGTase derived from *Bacillus circulans* Strain 251, which CGTase has been
complexed with acarbose, an effective CGTase inhibitor, and *Knegtel et al.* [*Knegtel R*
M A, Wind R D, Rozeboom H J, Kalk K H, Buitelaar R M, Dijkhuizen L and Dijkstra B
15 *W; J. Mol. Biol.* 1996 **256** 611-622] report the tertiary structure of a CGTase derived
from *Thermoanaerobacterium thermosulfurigenes.*

BRIEF DISCLOSURE OF THE INVENTION

The inventors have modified the amino acid sequence of a maltogenic alpha-
amylase to obtain variants with improved properties, based on the three-dimensional
20 structure of the maltogenic alpha-amylase Novamyl. The variants have altered
physicochemical properties., e.g. an altered pH optimum, improved thermostability,
increased specific activity, an altered cleavage pattern or an increased ability to reduce
retrogradation of starch or staling of bread.

Accordingly, the present invention provides a method of constructing a variant
25 of a parent maltogenic alpha-amylase, wherein the variant has at least one altered
property as compared to said parent maltogenic alpha-amylase, which method
comprises:

i) analyzing the structure of the maltogenic alpha-amylase to identify, on the
basis of an evaluation of structural considerations, at least one amino acid residue or at
30 least one structural region of the maltogenic alpha-amylase, which is of relevance for
altering said property;

ii) constructing a variant of the maltogenic alpha-amylase, which as compared
to the parent, has been modified in the amino acid residue or structural part identified in
i) so as to alter said property; and
35   iii) testing the resulting maltogenic alpha-amylase variant for said property.

The property which may be altered by the above methods of the present
invention may be, e.g., stability, pH dependent activity, ability to reduce retrogradation

of starch or staling of bread, specific activity, or substrate specificity. Thus, the variant may have, e.g., increased thermostability or higher activity at a lower pH an altered pH optimum, improved thermostability, increased specific activity or increased ability to reduce retrogradation of starch or staling of bread

5   In still further aspects the invention relates to variants of a maltogenic alpha-amylase, the DNA encoding such variants and methods of preparing the variants. Finally, the invention relates to the use of the variants for various industrial purposes, in particular baking.

DETAILED DISCLOSURE OF THE INVENTION

10 **Maltogenic alpha-amylase**

The maltogenic alpha-amylase is an enzyme classified in EC 3.2.1.133. The enzymatic activity does not require a non-reducing end on the substrate and the primary enzymatic activity results in the degradation of amylopectin and amylose to maltose and longer maltodextrins. It is able to hydrolyze amylose and amylopectin to

15 maltose in the alpha-configuration, and is also able to hydrolyze maltotriose as well as cyclodextrin.

A particularly preferred maltogenic alpha-amylase is the amylase cloned from *Bacillus* as described in EP 120 693 (hereinafter referred to as Novamyl). Novamyl has the amino acid sequence set forth in amino acids 1-686 of SEQ ID NO: 1. Novamyl is

20 encoded in the gene harbored in the *Bacillus* strain NCIB 11837 which has the nucleic acid sequence set forth in SEQ ID NO:1. The three-dimensional structure of Novamyl is described below.

In general, a preferred maltogenic alpha-amylase should have one or more of the following properties:

25   i) a three dimensional structural homology to Novamyl,

ii) an amino acid sequence having at least 70 % identity to SEQ ID NO: 1, preferably at least 80 % or 90 %, e.g. 95 % or 98 %,

iii) a DNA sequence which hybridizes to the DNA sequence set forth in SEQ ID NO:1 or to the DNA sequence encoding Novamyl harbored in the *Bacillus* strain NCIB

30 11837;

iv) a calcium binding site comprising a coordination equivalent to a backbone carbonyl atom from Asn77, sidechain atom OE2 and OE1 from Glu102, a sidechain atom OD1 from Asp79, a sidechain atom OD1 from Asp76, and a sidechain atom OE1 from Glu101, and one water molecule WAT V21, atom OW0, wherein the positions are

35 as shown in Appendix 1;

v) a sequence of five amino acids corresponding to Pro-Ala-Gly-Phe-Ser in a position equivalent to residues 191-195 in the amino acid sequence shown in SEQ ID NO: 1; and

The structural homology referred to above in i) is based on other sequence homologies, hydrophobic cluster analysis or by reverse threading (Huber, T ; Torda, AE, PROTEIN SCIENCE Vol. 7 , No. 1 pp. 142-149 (1998)) and which by any of these methods is predicted to have the same tertiary structure as Novamyl, wherein the tertiary structure refers to the overall folding or the folding of Domains A, B, and C, more preferably including Domain D, and most preferably including Domain E. Alternatively, a structural alignment between Novamyl and a maltogenic alpha-amylase may be used to identify equivalent positions.

The calcium binding site referred to above in iv) is based on a calcium binding site identified in the three-dimensional structure of Novamyl, and is discussed below in the section "Calcium binding sites."

The "equivalent position" referred to above in v) is based on amino acid or DNA sequence alignment or structural homology using methods known in the art.

Three-dimensional structure of maltogenic alpha-amylase

Novamyl was used to elucidate the three-dimensional structure forming the basis for the present invention.

The structure of Novamyl was solved in accordance with the principle for x-ray crystallographic methods, for example, as given in X-Ray Structure Determination, Stout, G.K. and Jensen, L.H., John Wiley & Sons, Inc. NY, 1989.

The structural coordinates for the solved crystal structure of Novamyl at 2.2 Å resolution using the isomorphous replacement method are given in standard PDB format (Protein Data Bank, Brookhaven National Laboratory, Brookhaven, CT) as set forth in Appendix 1. It is to be understood that Appendix 1 forms part of the present application. In the context of Appendix 1, the following abbreviations are used: CA refers to calcium ion or alpha-carbon atom of the polypeptide backbone, WAT refers to water or to calcium, MAL refers to maltose, HEX refers to a carbohydrate unit of a substrate analogue, and SUL refers to a sulfate ion.

Amino acid residues of the enzyme are identified herein by their respective one- or three-letter amino acid code.

The structure of said maltogenic alpha-amylase is made up of five globular domains, ordered A, B, C, D and E. The domains can be defined as being residues 1-132 and 204-403 for Domain A, residues 133-203 for Domain B, residues 404-496 for Domain C, residues 497-579 for Domain D, and residues 580-686 for Domain E,

wherein the numbering refers to the amino acid sequence in SEQ ID NO: 1. Features of Domains A, B, and C of particular interest are described below.

Domain A

Domain A is the largest domain and contains the active site which comprises a
5   cluster of three amino acid residues, D329, D228 and E256, spatially arranged at the bottom of a cleft in the surface of the enzyme. The structure of Domain A shows an overall fold in common with the α-amylases for which the structure is known, viz. the (beta/alpha) 8 barrel with eight central beta strands (numbered 1-8) and eight flanking a-helices. The β-barrel is defined by McGregor *op. cit.* The C-terminal end of the beta
10 strand 1 is connected to helix 1 by a loop denoted loop 1 and an identical pattern is found for the other loops, although the loops show some variation in size and some can be quite extensive.

The eight central beta-strands in the (beta/alpha) 8 barrel superimpose reasonably well with the known structures of CGTases. This part of the structure,
15 including the close surroundings of the active site located at the C-terminal end of the beta-strands, shows a high degree of identity with CGTases.

In contrast, the loops connecting the beta-strands and alpha helices display a high degree of variation from the known structures of CGTases. These loops constitute the structural context of the active site, and the majority of the contacts to the substrate
20 is found among residues located in these loops. Distinguishing characteristics such as substrate specificity, substrate binding, pH activity profile, substrate cleavage pattern, and the like, are determined by specific amino acids and the positions they occupy in these loops. In Novamyl Domain A contains two calcium binding sites, one of which is homologous to the calcium binding site in CGTases; the other is unique to Novamyl.
25 The structure of the calcium binding site is discussed further below in the section "Calcium binding sites."

Domain B

Domain B, also referred to as loop 3 of the (beta/alpha) 8 barrel, in comprises amino acid residues 133-203 of the amino acid sequence shown in SEQ ID NO: 1. The
30 structure is partially homologous to the structure of Domain B in CGTases, the most striking difference being the presence of a five amino acid insert corresponding to positions 191-195 in the amino acid sequence shown in SEQ ID NO: 1 which is not found in the CGTases. This insert is spatially positioned close to the active site residues and in close contact to the substrate.

Domain C

Domain C in Novamyl comprises amino acid residues 404-496 of the amino acid sequence shown in SEQ ID NO: 1. Domain C is composed entirely of β-strands which form a single 8-stranded sheet structure that folds back on itself, and thus may
5 be described as a β-sandwich structure. One part of the β-sheet forms the interface to Domain A.

Calcium binding sites

The structure of the maltogenic alpha-amylase exhibits three calcium-binding sites; that is, three calcium ions are found to be present in the structure. In common
10 with most of the known family 13 structures, one calcium ion, WAT 693 in Appendix 1, is located between the A and B domains. This calcium ion is coordinated by a backbone carbonyl atom from Gln184 and His232, sidechain atoms OD2 and OD1 from Asp198 , a sidechain atom OD1 from Asn131, and three water molecules WAT V1, WAT V5 and WAT V8.

15 A second calcium ion is located in the A domain and is common to CGTases, but not found in α-amylases. The calcium ion WAT 694 is coordinated by a backbone carbonyl atom from Gly48 and Asp23, sidechain atom OD2 from Asp50, a sidechain atom OD1 from Asp21, a sidechain atom OD1 from Asn26, and a sidechain atom OD1 from Asn27, and one water molecule WAT V62.

20 The third calcium ion is located in the A Domain and is unique to Novamyl. The calcium ion is WAT 692 and the coordination comprises a backbone carbonyl atom from Asn77, sidechain atom OE2 and OE1 from Glu102, a sidechain atom OD1 from Asp79, a sidechain atom OD1 from Asp76, and a sidechain atom OE1 from Glu101, and one water molecule WAT V21.

25 ## Substrate Binding Site

Parts of the loop discussed above in the context of domains A and B are of particular interest for substrate interaction and active site reactivity. In particular, in domain A, residues 37-45 in loop 1, residues 261-266 in loop 5, residues 327-330 in loop 7 and residues 370-376 in loop 8; in domain B, residues 135-145 in loop 3,
30 residues 173-180 and 188-196 in loop 3, wherein residue positions correspond to the amino acids in the amino acid sequence in SEQ ID NO: 1.

Without being limited to any theory, it is presently believed that binding between a substrate and an enzyme is supported by favorable interactions found within a sphere of 4 to 6 Å between the substrate molecule and the enzyme, such as
35 hydrogen bonds and/or strong electrostatic interaction. The following residues of

Novamyl (SEQ ID NO: 1), are within a distance of 6 Å of the substrate HEX and thus believed to be involved in interactions with said substrate:

      44, 89, 90, 92, 93, 127, 129, 132, 135, 177, 178, 188, 191, 194, 196, 226, 228, 229, 230, 231, 232, 256, 258-261, 288, 328, 329, 371, 372, 373, 376, and 690.

      The following residues of Novamyl are within a distance of 4 Å of the substrate HEX and thus believed to be involved in interactions with said substrate:

      90, 92, 93, 129, 132, 177, 188, 189, 190, 191, 196, 226, 228, 229, 231, 232, 256, 258, 259, 260, 261, 328, 329, 372, 376, and 690.

Homology building of Novamyl®

      The structure of the Novamyl® was model built on the structure disclosed in Appendix 1 herein. The structure of other maltogenic alpha-amylases may be built analogously.

      A model structure of a maltogenic alpha-amylase can be built using the Homology program or a comparable program, eg., Modeller (both from Molecular Simulations, Inc., San Diego, CA). The principle is to align the sequence of the maltogenic alpha-amylase with the known structure with that of the maltogenic alpha-amylase for which a model structure is to be constructed. The structurally conserved regions can then be built on the basis of consensus sequences. In areas lacking homology, loop structures can be inserted, or sequences can be deleted with subsequent bonding of the necessary residues using, e.g., the program Homology. Subsequent relaxing and optimization of the structure should be done using either Homology or another molecular simulation program, e.g., CHARMm from Molecular Simulations.

Methods for designing novel maltogenic alpha-amylase variants

      In a first aspect, the invention relates to a method of constructing a variant of a parent maltogenic alpha-amylase, wherein said variant has at least one altered property as compared to said parent α-amylase, which method comprises:

      i) analyzing the structure of the maltogenic alpha-amylase to identify at least one amino acid or structural region of said α-amylase, which, on the basis of structural or functional considerations, is determined to be of relevance for altering said property of the parent maltogenic alpha-amylase;

      ii) constructing a variant of the maltogenic alpha-amylase, which as compared to the parent, has been modified in the amino acid residue or structural region identified in i) has been modified so as to alter said property; and

      iii) testing the resulting variant for said property.

The structural part which is identified in step i) of the method of the invention may be composed of one amino acid residue. However, normally the structural part comprises more than one amino acid residue, typically constituting one of the above parts of the maltogenic alpha-amylase structure such as one of the A, B, C, D or E

5   domains, an interface between any of these domains, a calcium binding site, a loop structure, the substrate binding site, or the like.

The structural or functional considerations may involve an analysis of the relevant structure or structural part and its contemplated impact on the function of the enzyme. For example, an analysis of the functional differences between maltogenic

10 alpha-amylase and the various CGTases may be used for assigning certain properties of Novamyl to certain parts of the Novamyl structure or to contemplate such relationship. For instance, differences in the pattern or structure of loops surrounding the active site may result in differences in access to the active site of the substrate and thus differences in substrate specificity and/or cleavage pattern.

15 Furthermore, parts of a maltogenic alpha-amylase involved in substrate binding, and thus, for example, substrate specificity and/or cleavage, calcium ion binding, important, for example, for the calcium dependency of the enzyme, and the like, have been identified (vide infra).

The modification of an amino acid residue or structural region is typically

20 accomplished by suitable modifications of a DNA sequence encoding the parent enzyme in question. The modification may be substitution, deletion or insertion of an amino acid residue or a structural part.

The property to be modified may be stability (e.g. thermostability), pH dependent activity, substrate specificity, specific activity or ability to reduce

25 retrogradation of starch or staling of bread. Thus, the altered property may be an altered specific activity at a given pH and/or an altered substrate specificity, such as an altered pattern of substrate cleavage or an altered pattern of substrate inhibition.

In step ii) of the method according to the invention the part of the structure to be identified is preferably one which in the folded enzyme is believed to be in contact

30 with the substrate (cf, the disclosure above in the section entitled "Substrate Binding Site") or involved in substrate specificity and/or cleavage pattern, and/or one which is in contact with one of the calcium ions and/or one, which is contributing to the pH or temperature profile of the enzyme, or is otherwise responsible for the properties of the maltogenic alpha-amylase.

35 Described in the following are specific types of variants which have been designed by use of the method of the invention.

The variants of the invention may comprise additional modifications in addition to the modifications described herein. The variants preferably have an amino acid

having more than 70 % identity with SEQ ID NO: 1, preferably more than 80 %, particularly more than 90 %, especially more than 95 %, e.g. more than 98 %.

Maltogenic alpha-amylase variants with altered pH dependent activity profile

The pH dependent activity profile can be changed by changing the pKa of

5 residues within 10 Å of the active site residues of the maltogenic alpha-amylase. Changing the pKa of the active site residues is achieved, e.g., by changing the electrostatic interaction or hydrophobic interaction between functional groups of amino acid side chains of a given amino acid residue and its close surroundings. To obtain a higher activity at a higher pH, negatively charged residues are placed near a hydrogen

10 donor acid, whereas positively charged residues placed near a nucleophilic acid will result in higher activity at low pH. Also, a decrease in the pKa can be obtained by reducing the accessibility of water or increasing hydrophobicity of the environment.

Thus, another aspect of the present invention relates to a variant of a parent maltogenic alpha-amylase, in which the variant has an altered pH dependent activity

15 profile as compared to the parent, wherein the variant may be obtained by the following method:

i) identifying an amino acid residue within 15 Å from an active site residue of a maltogenic alpha-amylase in the three-dimensional structure of said parent maltogenic alpha-amylase, in particular 10 Å from an active site residue, wherein said amino acid

20 residue is contemplated to be involved in electrostatic or hydrophobic interactions with an active site residue;

ii) substituting, in the structure, said amino acid residue with an amino acid residue which changes the electrostatic and/or hydrophobic surroundings of an active site residue, and evaluating the accommodation of the amino acid residue in the

25 structure,

iii) optionally repeating step i) and/or ii) recursively until an amino acid substitution has been identified which is accommodated into the structure,

iv) constructing a maltogenic alpha-amylase variant resulting from steps i) and ii), and optionally iii), and testing the pH dependent enzymatic activity of said variant.

30 In a preferred embodiment, the variant of a maltogenic alpha-amylase having an altered pH dependent activity profile as compared to the parent maltogenic alpha-amylase comprises a modification of an amino acid residue corresponding to one or more of the following residues of the amino acid sequence set forth in SEQ ID NO: 1:

D127, V129, F188, A229, Y258, V281, F284, T288, N327, M330, G370, N371,

35 and D372,

L71, S72, V74, L75, L78, T80, L81, G83, T84, D85, N86, T87, G88, Y89, H90, G91, T94, R95, D96, F97, I174, S175, N176, D178, D179, R180, Y181, E182, A183, Q184, K186, N187, F188, T189, D190, A192, G193, F194, S195, L196.

In more preferred embodiment, the variant comprises a modification
5 corresponding to one or more of the following modifications in the amino acid sequence set forth in SEQ ID NO: 1:

D127N/L, V129S/T/G/V, F188E/K/H, A229S/T/G/V, Y258E/D/K/R/F/N, V281L/T, F284K/H/D/E/Y, T288E/K/R, N327D, M330L/F/I/D/E/K, G370N, N371D/E/G/K, and D372N/V,

10 L71I, S72C, V74I, L75N/D/Q/I/V, L78N/I, T80I/L/V/S/N/G, L81I/V/S/T/N/Q/K/H, G83A/S/T/N/Q/E/D/R/H/L, T84S/A/N/D/G, D85A/T/S/N/G, N86Q/E/D/Y/H/K, T87S/I, G88A/S/T, Y89F, H90N/Q/K, G91A/S/T, T94N/D/A/M/V/I, R95K/Q, D96N/V/Q/I, F97Y, I174N/Q/L, S175T/A/N/D, N176S/T/H/Q/P, D178N/Q/E/K/H, D179Y/N/H, R180W, Y181R/F/C/L, E182D, A183S/C/G, Q184E, K186R, N187Q/E/L/F/H/K/V/L,
15 F188Y/L/I/H/N, T189N/D/A/S/H/Y/G, D190E/Q/H/N/K, A192T/D/E/N/K, G193A/S/T, F194Y, S195N/D/E/R/K/G, L196I.

Similar modifications may be introduced in equivalent positions of other maltogenic alpha-amylases. Variants of particular interest have a combination of one or more of the above with any of the other modifications disclosed herein.

20 **Maltogenic alpha-amylase variants with altered stability**

A variant with improved stability (typically increased stability) may be obtained by stabilization of calcium binding, substitution with proline, substitution of histidine with another amino acid, introduction of an interdomain disulfide bond, removal of a deamidation site, altering a hydrogen bond contact, filling in an internal structural cavity
25 with one or more amino acids with bulkier side groups, introduction of interdomain interactions, altering charge distribution, helix capping, or introduction of a salt bridge.

Calcium binding

The invention provides a variant of a parent maltogenic alpha-amylase, which has an altered stability due to an altered stabilization of calcium (Ca^{2+}) binding. The
30 enzyme variant may have altered thermostability or pH dependent stability, or it may have maltogenic alpha-amylase activity in the presence of a lower concentration of calcium ion. It is presently believed that amino acid residues located within 10 Å from a calcium ion are involved in or are of importance for the Ca^{2+} binding capability of the enzyme.

The amino acid residues found within a distance of 10 Å from the Ca^{2+} binding sites of the maltogenic alpha-amylase with the amino acid sequence set forth in SEQ ID NO: 1 were determined as described in Example 2 and are as follows:

16, 17, 18, 19, 20, 21, 22, 23, 24, 25, 26, 27, 28 29, 30, 31, 32, 33,35, 36, 40, 46, 47, 48, 49, 50, 51, 52, 53, 54, 56, 73, 74, 75, 76, 77, 78, 79, 80, 81, 87, 88, 89, 91, 93, 94, 95, 96, 99, 100, 101, 102, 103, 104, 105, 109, 129, 130, 131, 132, 133, 134, 145, 150, 167, 168, 169, 170, 171, 172, 174, 177, 180, 181, 182, 183, 184, 185, 186, 187, 188, 189, 196, 197, 198, 199, 200, 201, 202, 206, 210, 228, 229, 230, 231, 232, 233, 234, 235, 237, 378, and 637.

In order to construct a variant according to this aspect of the invention it is desirable to substitute at least one of the above mentioned amino acid residues, which is determined to be involved in a non-optimal calcium binding, with any other amino acid residue which improves the Ca^{2+} binding affinity of the variant enzyme. Accordingly, another aspect of the invention relates to a method of constructing a variant of a parent maltogenic alpha-amylase wherein said variant has a stabilised Ca^{2+} binding as compared to said parent, which method comprises:

i) identifying an amino acid residue within 10 Å from a Ca^{2+} binding site of a maltogenic alpha-amylase in a model of the three-dimensional structure of said α-amylase which, from structural or functional considerations, is determined to be responsible for a non-optimal calcium ion interaction;

ii) constructing a variant in which said amino acid residue is substituted with another amino acid residue which, from structural or functional considerations, is determined to be important for establishing an altered Ca^{2+} binding affinity; and

iii) testing the Ca^{2+} binding of the resulting maltogenic alpha-amylase variant.

Substituting an amino acid residue responsible for non-optimal calcium ion interaction with another residue may alter a calcium ion binding interaction of the enzyme. For instance, the amino acid residue in question may be selected on the basis of one or more of the following objectives:

a) to obtain an improved interaction between a calcium ion and an amino acid residue as identified from the structure of the maltogenic alpha-amylase. For instance, if the amino acid residue in question is exposed to a surrounding solvent, it may be advantageous to increase the shielding of said amino acid residue from the solvent so as to stabilize the interaction between said amino acid residue and a calcium ion. This can be achieved by substituting said residue, or an amino acid residue in the vicinity of said residue contributing to the shielding, with an amino acid residue with a bulkier side group or which otherwise results in an improved shielding effect.

b) to stabilize a calcium binding site, for instance by stabilizing the structure of the maltogenic alpha-amylase, e.g. by stabilizing the contacts between two or more of

the five domains or stabilizing one or more of the individual domains as such. This may, e.g., be achieved by providing for a better coordination to amino acid side chains, which may, e.g., be obtained by substituting an N residue with a D residue and/or a Q residue with an E residue, e.g. within 10 Å, and preferably within 3 or 4 Å, of a calcium binding 5   site.

c) to improve the coordination between the calcium ion and the calcium binding residues, e.g., by improving the interaction between the ion and the coordinating residues or increasing the number of sidechain coordinations by substituting a coordinating water with an amino acid sidechain.

10   d) replace water by a coordinating calcium amino acid residue.

Preferably, the amino acid residue to be modified is located within 8 Å of a Ca^{2+} ion, preferably within 5 Å of a Ca^{2+} ion. The amino acid residues within 8 Å and 5 Å, respectively, may easily be identified by an analogous method used for identifying amino acid residues within 10 Å (cf. Example 2).

15   In a preferred embodiment, the variant of a maltogenic alpha-amylase having an altered Ca^{2+} binding as compared to the parent maltogenic alpha-amylase comprises a substitution of an amino acid residue corresponding to one or more of the following residues of the amino acid sequence set forth in SEQ ID NO: 1:

D17, A30, S32, R95, H103, N131, Q201, I174, and/or H169,

20   V74, L75, L78, T80, L81, T87, G88, Y89, H90, G91, T94, R95, D96, F97, Y167, F168, H169, H170, N171, G172, D173, I174, S175, N176, D178, D179, R180, Y181, E182, A183, Q184, K186, N187, F188, T189.

In more preferred embodiment, the variant of a maltogenic alpha-amylase comprises a substitution corresponding to one or more of the following substitutions in 25   the amino acid sequence set forth in SEQ ID NO: 1:

D17E/Q, A30M/L/A/V/I/E/Q, S32D/E/N/Q, R95M/L/A/V/I/E/Q, H103Y/N/Q/D/E, N131D, Q201E, I174E/Q, and H169N/D/E/Q

V74I, L75N/D/Q/I/V, L78N/I, T80I/L/V/S/N/G, L81I/V/S/T/N/Q/K/H, T87S/I, G88A/S/T, Y89F, H90N/Q/K, G91A/S/T, T94N/D/A/M/V/I, R95K/Q, D96N/V/Q/I, F97Y, 30   Y167F/R/C, F168Y, H169N/Q/K, H170N/Q/K, N171D/E/Q/H/R/K/G, G172A/T/S, D173N/S/T/Y/R/G, I174N/Q/L, S175T/A/N/D, N176S/T/H/Q/P, D178N/Q/E/K/H, D179Y/N/H, R180W, Y181R/F/C/L, E182D, A183S/C/G, Q184E, K186R, N187Q/E/L/F/H/K/V/L, F188Y/L/I/H/N, T189N/D/A/S/H/Y/G.

In another preferred embodiment of the invention with respect to altering the 35 Ca^{2+} binding of a maltogenic alpha-amylase the partial sequence N28-P29-A30-K31-S32-Y33-G34 as set forth in SEQ ID NO: 1 is modified.

Similar substitutions may be introduced in equivalent positions of other maltogenic alpha-amylases. Modifications of particular interest are any combination of one or more of the above with any of the other modifications disclosed herein.

Other substitutions

5   Variants with improved stability of the enzyme can be achieved by improving existing or introducing new interdomain and intradomain contacts. Such improved stability can be achieved by the modifications listed below.

The maltogenic alpha-amylase having the amino acid sequence shown in SEQ ID NO: 1 may be stabilized by the introduction of one or more interdomain disulfide
10 bonds. Accordingly, another preferred embodiment of the present invention relates to a variant of a parent maltogenic alpha-amylase which has improved stability and at least one more interdomain disulfide bridge as compared to said parent, wherein said variant comprises a modification in a position corresponding to at least one of the following pairs of positions in SEQ ID NO: 1:

15   G236 + S583, G618 + R272, T252 + V433 and/or A348 + V487.

In a more preferred embodiment, the substitution corresponds to at least one of the following pairs:

G236C + S583C, G618C + R272C, T252C + V433C and/or A348C + V487C.

Another preferred embodiment of the invention relates to a variant of a parent
20 maltogenic alpha-amylase which has an improved stability and an altered interdomain interaction as compared to said parent, wherein said variant comprises a substitution in a position corresponding at least one of the following sets of positions in SEQ ID NO: 1:

i) F143, F194, L78;
ii) A341, A348, L398, I415, T439, L464, L465;
25   iii) L557;
iv) S240, L268;
v) Q208, L628;
vi) F427, Q500, N507, M508, S573; and
vii) I510, V620.

30   In a more preferred embodiment, the substitution corresponds to at least one of the following sets:

i) F143Y, F194Y, L78Y/F/W/E/Q;
ii) A341S/D/N, A348V/I/L, L398E/Q/N/D, I415E/Q, T439D/E/Q/N, L464D/E, L465D/E/N/Q/R/K;
35   iii) L557Q/E/N/D;
iv) S240D/E/N/Q, L268D/E/N/Q/R/K;
v) Q208D/E/Q, L628E/Q/N/D;

vi) F427E/Q/R/K/Y, Q500Y, N507Q/E/D, M508K/R/E/Q, S573D/E/N/Q; and/or
vii) I510D/E/N/Q/S, V620D/E/N/Q.

Another preferred embodiment of the invention relates to a variant of a parent maltogenic alpha-amylase which has an improved stability and one or more salt bridges as compared to said parent, wherein said variant comprises a substitution in a position corresponding at least one of the following sets of positions in SEQ ID NO: 1:

N106, N320 and Q624.

In a more preferred embodiment, the variant of a maltogenic alpha-amylase comprises a substitution corresponding to the following substitutions in the amino acid sequence set forth in SEQ ID NO: 1:

N106R, N320E/D and/or Q624E.

Another embodiment of the invention relates to a variant of a parent maltogenic alpha-amylase which has an improved stability and wherein said variant comprises a substitution in a position corresponding at least one of the following sets of positions in SEQ ID NO: 1:

K40, V74, S141, T142, F188, N234, K249, D261, D261, L268, V279, N342, G397, A403, K425, S442, S479, S493, T494, S495, A496, S497, A498, Q500, K520, A555 and N595.

In a more preferred embodiment, the variant of a maltogenic alpha-amylase comprises a substitution corresponding to one or more of the following substitutions with proline in the amino acid sequence set forth in SEQ ID NO: 1:

V74P, S141P, N234P, K249P, L268P, V279P, N342P, G397P, A403P, S442P, S479P, S493P, T494P, S495P, A496P, S497P, A498P, Q500P, and/or A555P.

Other preferred substitutions are K40R, T142A, F188I/L, D261G, K425E, K520R, and/or N595I.

Analogously, it may be preferred that one or more histidine residues present in the parent maltogenic alpha-amylase is or are substituted with a non-histidine residues such as Y, V I, L, F, M, E, Q, N, or D. Accordingly, in another preferred embodiment, the variant of a maltogenic alpha-amylase comprises a substitution of an amino acid residue corresponding to one or more of the following residues of the amino acid sequence set forth in SEQ ID NO: 1:

H103, H220, and H344

In a more preferred embodiment, the variant of a maltogenic alpha-amylase comprises a substitution corresponding to one or more of the following substitutions in the amino acid sequence set forth in SEQ ID NO: 1:

H103Y/V/I/L/F/Y, H220Y/L/M, and H344E/Q/N/D/Y.

It may be preferred that one or more asparagine or glutamine residues present in the parent maltogenic alpha-amylase is or are substituted with a residue lacking the

amide on the side chain. Accordingly, in another preferred embodiment, the variant of a Novamyl-like comprises a substitution of an amino acid residue corresponding to one or more of the following residues of the amino acid sequence set forth in SEQ ID NO: 1:

Q13, N26, N77, N86, N99, Q119, N120, N131, N152, N171, N176, N187,
5   Q201, N203, N234, Q247, N266, N275, N276, N280, N287, Q299, N320, N327, N342, Q365, N371, N375, N401, N436, N454, N468, N474, Q500, N507, N513, Q526, N575, Q581, N621, Q624 and N664.

In more preferred embodiment, the variant of a maltogenic alpha-amylase comprises a substitution corresponding to one or more of the following substitutions in
10 the amino acid sequence set forth in SEQ ID NO: 1:

Q13S/T/A/V/L/I/F/M,   N26S/T/A/V/L/I,   N77S/T/A/V/L/I,   N86S/T/A/V/L/I,
N99T/S/V/L, Q119T/S, N120S/T/A/V/L/I, N131S/T/A/V/L/I, N152T/S/V/L, N171Y/D/S/T, N176S/T/A/V/L/I,   N187S/T/A/V/L/I,   Q201S/T/A/V/L/I/F/M,   N203D/S/T/A/V/L/I, N234S/T/A/V/L/I,   Q247S/T/A/V/L/I/F/M,   N266S/T/A/V/L/I,   N275S/T/A/V/L/I,
15 N276S/T/A/V/L/I, N280S/T/A/V/L/I, N287S/T/A/V/L/I, Q299L/T/S, N320S/T/A/V/L/I, N327S/T/A/V/L/I, N342S/T/A/V/L/I, Q365S/T/A/V/L/I, N371S/T/A/V/L/I, N375S/T/A/V/L/I, N401S/T/A/V/L/I,   N436S/T/A/V/L/I,   N454D/S/T/A/V/L/I,   N468D/S/T/A/V/L/I, N474D/S/T/A/V/L/I, Q500S/T/A/V/L/I/F/M, N507S/T/A/V/L/I, N513S/T/A/V/L/I, Q526 D/S/T/A/V/L/I,   N575S/T/A/V/L/I,   Q581S/T/A/V/L/I/F/M,   N621S/T/A/V/L/I
20 Q624S/T/A/V/L/I/F/M and N664D/S/T/A/V/L/I.

Another embodiment of the invention relates to a variant of a parent maltogenic alpha-amylase which has improved stability and improved hydrogen bond contacts as compared to said parent, wherein said variant comprises a modification in a position corresponding to one or more of the following positions in SEQ ID NO: 1:
25   I16, L35, M45, P73, D76, D79, A192, I100, A148, A163+G172, L268, V281, D285, L321, F297, N305, K316, S573, A341, M378, A381, F389, A483, A486, I510, A564, F586, K589, F636, K645, A629, and/or T681.

In a preferred embodiment, the modification corresponds to one or more of the following:
30   I16T/D/N, L35Q, M45K, P73Q, D76E, D79E/Y, A192S/D/N, I100T/S/D/N/E/Q, A148D/N/E/Q/S/T/R/K, A163Y+G172S/D/N, L268R/K, V281/Q, D285R/K, L321Q, F297N/D/Q/E, N305K/R, K316N/D, S573N/D, A341R/K, M378R/K, A381S/D/N, F389Y, A483S/D/N, A486Q/E, I510R/K, A564S/D/N, F586S/D/N, K589S/D/Q/N, F636Y, K645T, A629N/D/E/Q, and/or T681D/N/E/Q/S.

35   Similar substitutions may be introduced in equivalent positions of other maltogenic alpha-amylases. Substitutions of particular interest are any combination of one or more of the above with any of the other modifications disclosed herein.

Before actually constructing a maltogenic alpha-amylase variant to achieve any of the above objectives, it may be convenient to evaluate whether or not the contemplated amino acid modification can be accommodated into the maltogenic alpha-amylase structure, e.g. into a model of the three-dimensional structure of the parent maltogenic alpha-amylase.

Maltogenic alpha-amylase variants with altered thermostability and/or altered temperature dependent activity profile

The invention further relates to a variant of a parent maltogenic alpha-amylase, which results from substitution, deletion or insertion of one or more amino acid residues so as to obtain a variant having an altered thermostability or temperature dependent activity profile.

The structure of the maltogenic alpha-amylase contains a number of unique internal cavities which may contain water and a number of crevices. In order to increase the thermostability of the polypeptide it may be desirable to reduce the number or size of cavities and crevices, e.g., by introducing one or more hydrophobic contacts, preferably achieved by introducing amino acids with bulkier side groups in the vicinity or surroundings of the cavity. For instance, the amino acid residues to be modified are those which are involved in the formation of the cavity.

Accordingly, in a further aspect the present invention relates to a method of increasing the thermostability and/or altering the temperature dependent activity profile of a parent maltogenic alpha-amylase, which method comprises:

i) identifying an internal cavity or a crevice of the parent maltogenic alpha-amylase in the three-dimensional structure of said polypeptide;

ii) substituting, in the structure, one or more amino acid residues in the neighbourhood of the cavity or crevice identified in step i) with another amino acid residue which, from structural or functional considerations, is determined to increase the hydrophobic interaction and to fill out or reduce the size of the cavity or crevice; and

iii) constructing a variant of the parent maltogenic alpha-amylase resulting from step ii) and testing the thermostability and/or temperature dependent activity of the variant.

The structure identified in Appendix 1 may be used for identifying the cavity or crevice of the parent maltogenic alpha-amylase.

It will be understood that the cavity or crevice is identified by the amino acid residues surrounding said cavity or crevice, and that modification of said amino acid residues are of importance for filling or reducing the size of said cavity or crevice. Preferably, the modification is a substitution with a bulkier amino acid residue, i.e. one with a greater side chain volume. For example, all the amino acids are bulkier than Gly,

whereas Tyr and Trp are bulkier than Phe. The particular amino acid residues referred to below are those which in a crystal structure have been found to flank the cavity or crevice in question.

In a preferred embodiment, the variant of a maltogenic alpha-amylase, in order
5   to fill, either completely or partly, cavities located internally in the structure, comprises a substitution of an amino acid residue corresponding to one or more of the following residues of the amino acid sequence set forth in SEQ ID NO: 1:

L51, L75, L78, G88, G91, T94, V114, I125, V126, T134, G157, L217, S235, G236, V254, V279, V281, L286, V289, I290, V308, L321, I325, D326, L343, F349,
10 S353, I359, I405, L448, Q449, L452, I470, G509, V515, S583, G625, L627, L628 and A670.

L71, S72, V74, L75, L78, T80, L81, G83, T84, D85, N86, T87, G88, Y89, H90, G91, T94, R95, D96, F97, Y167, F168, H169, H170, N171, G172, D173, I174, S175, N176, D178, D179, R180, Y181, E182, A183, Q184, K186, N187, F188, T189, D190,
15 A192, G193, F194, S195, L196.

In a more preferred embodiment, the variant of a maltogenic alpha-amylase comprises one or more substitutions corresponding to the following substitutions in the amino acid sequence set forth in SEQ ID NO: 1:

L217 in combination with L75 (e.g. L217F/Y in combination with L75F/Y),
20 L51W, L75F/Y, L78I, G88A/V/T, G91T/S/V/N, T94V/I/L, V114V/I/L, I125L/M/F/Y/W, V126I/L, T134V/I/L/M/F/Y/W, G157A/V/I/L, L217V/I/M/F/Y/W, S235I/L/M/F/Y/W, G236A/V/I/L/M/F/Y/W, V254I/L/M/F/Y/W, V279M/I/L/F, V281I/L/M/F/Y/W, L286F, V289I/L/R, I290M/L/F, V308I/L/M/F/Y/W, L321I/M/F/Y/W, I325L/M/F/Y/W, D326E/Q, L343M/F/Y/W, F349W/Y, S353V/I/L, I359L/M/F/Y/W, I405M/L/Y/F/W, L448Y, Q449Y,
25 L452M/Y/F/W, I470M/L/F, G509A/V/I/L/M/S/T/D/N, V515I/L, S583V/I/L/V, G625A/V/I/L/M/F/Y/W, L627M/F/Y, L628M/I/F/Y/W and A670V/I/L/M/F/Y/W,

L71I, S72C, V74I, L75N/D/Q/I/V, L78N/I, T80I/L/V/S/N/G, L81I/V/S/T/N/Q/K/H, G83A/S/T/N/Q/E/D/R/H/L, T84S/A/N/D/G, D85A/T/S/N/G, N86Q/E/D/Y/H/K, T87S/I, G88A/S/T, Y89F, H90N/Q/K, G91A/S/T, T94N/D/A/M/V/I, R95K/Q, D96N/V/Q/I, F97Y,
30 Y167F/R/C, F168Y, H169N/Q/K, H170N/Q/K, N171D/E/Q/H/R/K/G, G172A/T/S, D173N/S/T/Y/R/G, I174N/Q/L, S175T/A/N/D, N176S/T/H/Q/P, D178N/Q/E/K/H, D179Y/N/H, R180W, Y181R/F/C/L, E182D, A183S/C/G, Q184E, K186R, N187Q/E/L/F/H/K/V/L, F188Y/L/I/H/N, T189N/D/A/S/H/Y/G, D190E/Q/H/N/K, A192T/D/E/N/K, G193A/S/T, F194Y, S195N/D/E/R/K/G, L196I.

35 Similar substitutions may be introduced in equivalent positions of other maltogenic alpha-amylases. Variants of particular interest have a combination of one or more of the above with any of the other modifications disclosed herein.

Maltogenic alpha-amylase variants with an alter d cleavag pattern

One aim of the present invention is to change the degradation characteristics of a maltogenic alpha-amylase. Thus, Novamyl hydrolyzes starch to form predominantly maltose (G2) and a small amount of glucose (G1), but virtually no higher
5 oligosaccharides (G3+). It may be desirable to change this cleavage pattern, e.g. so as to form higher amounts of higher oligosaccharides, such as maltotriose (G3), maltotetraose (G4) and maltopentaose (G5).

A variant of a parent maltogenic alpha-amylase in which the substrate cleavage pattern is altered as compared to said parent may be constructed by a method which
10 comprises:

i) identifying the substrate binding area of the parent maltogenic alpha-amylase in a model of the three-dimensional structure, e.g. within a sphere of 4 Å from the substrate binding site as defined in the section above entitled "Substrate Binding Site";

ii) substituting in the model one or more amino acid residues of the substrate
15 binding area of the cleft identified in i) which is or are believed to be responsible for the cleavage pattern of the parent with another amino acid residue which from structural or functional considerations is believed to result in an altered substrate cleavage pattern, or deleting one or more amino acid residues of the substrate binding area contemplated to introduce favorable interactions to the substrate or adding one or more
20 amino acid residues to the substrate binding area contemplated to introduce favorable interactions to the substrate; and

iii) constructing a maltogenic alpha-amylase variant resulting from step ii) and testing the substrate cleavage pattern of the variant.

Accordingly, another aspect of the invention relates to a variant of a parent
25 maltogenic alpha-amylase which has an altered substrate binding site as compared to said parent, which variant comprises a modification in a position corresponding to one or both of the following positions in SEQ ID NO: 1:

V281 and/or A629.

In a preferred embodiment, the variant comprises a modification corresponding
30 to:

V281Q and/or A629N/D/E/Q.

Similar modifications may be introduced in equivalent positions of other maltogenic alpha-amylases. Substitutions of particular interest are any combination of one or both of the above with any of the other modifications disclosed herein.

**Maltogenic alpha-amylase variants with improved ability to reduc retrograda-
tion of starch and/or staling of bread**

The invention provides maltogenic alpha-amylase variants having improved
ability to reduce the retrogradation of starch and/or the staling of bread. Preferred
5 variants comprise a modification at one or more positions corresponding to the
following amino acid residues in SEQ ID NO: 1:

A30, K40, N115, T142, F188, T189, P191, A192, G193, F194, S195, D261,
N327, K425, K520 and N595.

In a more preferred embodiment, the variant comprises one or more
10 modifications corresponding to the following in SEQ ID NO: 1:

A30D, K40R, N115D, T142A, F188L, T189Y, Δ (191-195), D261G, D261G,
N327S, K425E, K520R and N595I.

Determination of residues within 10Å from calcium ions

The coordinates of Appendix 1 are read into the INSIGHT program (BIOSYM
15 Technologies). The spatial coordinates are presented showing the bonds between the
atoms. The ions are presented as well as the water atoms. The part of the program
package for creating subsets was used to create a 10 Å subset around the calcium ions
in the structure by using the command ZONE. All residues identified as having an atom
within the designated 10 Å distance from a calcium ion are compiled and listed by using
20 the command LIST MOLECULE. By giving the ions the name "VAT CA" in the
coordinate file, a 10 Å sphere around all atoms called "VAT CA" is compiled. The
specific residues identified in this manner are given further above in the section entitled
"Calcium binding".

25 **Determination of cavities**

The solved structure of Novamyl with the structural coordinates set forth in
Appendix 1 reveals many internal crevices and cavities. When analysing for such
cavities the Connolly program is normally used (Lee, B. and Richards, F.M. (1971) J.
Mol. Biol. 55:379-400). The program uses a probe with radius to search the external
30 and internal surface of the protein. The smallest crevice observable in this way has the
probe radius.

To analyse the solved structure a modified version of the Connolly program
included in the program of INSIGHT was used. In the first step, the water molecules
and the ions were removed by unmerging these atoms from the solved structure. By
35 using the command MOLECULE SURFACE SOLVENT the solvent accessible surface
area was calculated for all atoms and residues using a probe radius of 1.4 Å, and

displayed graphically together with the model of the solved structure. The internal cavities are then seen as dot surfaces with no connections to the external surface.

Suggestions for specific modifications to fill out the cavities are given above in the section entitled "Variants with altered thermostability and/or altered temperature
5 dependent activity profile"). By using the homology built structures or/and comparisons based on sequence alignment, mutations for homologous structures of maltogenic alpha-amylases can be made.

Nomenclature for amino acid modifications

10 The nomenclature used herein for defining mutations is essentially as described in WO 92/05249. Thus, F188H indicates a substitution of the amino acid F (Phe) in position 188 with the amino acid H (His). V129S/T/G/V indicates a substitution of V129 with S, T, G or V. Δ (191-195) or Δ (191-195) indicates a deletion of amino acids in positions 191-195. 192-A-193 indicates an insertion of A between amino acids
15 192 and 193.

Polypeptide sequence identity

For purposes of the present invention, the degree of identity may be suitably determined according to the method described in Needleman, S.B. and Wunsch, C.D., (1970), Journal of Molecular Biology, 48, 443-45, with the following settings for
20 polypeptide sequence comparison: GAP creation penalty of 3.0 and GAP extension penalty of 0.1. The determination may be done by means of a computer program known such as GAP provided in the GCG program package (Program Manual for the Wisconsin Package, Version 8, August 1994, Genetics Computer Group, 575 Science Drive, Madison, Wisconsin, USA 53711).

25 The variants of the invention have an amino acid identity with amino acids 1-686 of SEQ ID NO: 1 of at least 70 %, preferably at least 80 %, e.g. at least 90 %, particularly at least 95 % or at least 98 %.

Hybridization

Suitable experimental conditions for determining hybridization between a
30 nucleotide probe and a homologous DNA or RNA sequence involves presoaking of the filter containing the DNA fragments or RNA to hybridize in 5x SSC (sodium chloride/sodium citrate, Sambrook, et al., 1989) for 10 min, and prehybridization of the filter in a solution of 5x SSC, 5x Denhardt's solution (Sambrook, et al., 1989), 0.5% SDS and 100 μg/ml of denatured sonicated salmon sperm DNA (Sambrook, et al.,
35 1989), followed by hybridization in the same solution containing a random-primed

(Feinberg, A. P. and Vogelstein, B. (1983) *Anal. Biochem.* 132:6-13), ^{32}P-dCTP-labeled (specific activity > 1 x 10^9 cpm/μg) probe for 12 hours at ca. 45°C. The filter is then washed twice for 30 minutes in 2x SSC, 0.5% SDS at least 55°C (low stringency), preferably at least 60°C (medium stringency), more preferably at least 65°C
5 (medium/high stringency), more preferably at least 70°C (high stringency), even more preferably at least 75°C (very high stringency).

Molecules which hybridize to the oligonucleotide probe under these conditions are detected by exposure to x-ray film.

Methods of preparing variants of maltogenic alpha-amylases

10 Cloning a DNA sequence encoding a Novamyl-like polypeptide

The DNA sequence encoding a parent maltogenic alpha-amylase may be isolated from any cell or microorganism producing the maltogenic alpha-amylase in question, using various methods well known in the art, for example, from the *Bacillus* strain NCIB 11837.

15 First, a genomic DNA and/or cDNA library should be constructed using chromosomal DNA or messenger RNA from the organism that produces the maltogenic alpha-amylase to be studied. Then, if the amino acid sequence of the α-amylase is known, homologous, labelled oligonucleotide probes may be synthesised and used to identify maltogenic alpha-amylase-encoding clones from a genomic library prepared
20 from the organism in question. Alternatively, a labelled oligonucleotide probe containing sequences homologous to a known α-amylase gene could be used as a probe to identify maltogenic alpha-amylase-encoding clones, using hybridization and washing conditions of lower stringency.

Another method for identifying maltogenic alpha-amylase-encoding clones
25 involves inserting fragments of genomic DNA into an expression vector, such as a plasmid, transforming α-amylase negative bacteria with the resulting genomic DNA library, and then plating the transformed bacteria onto agar containing a substrate for maltogenic alpha-amylase, thereby allowing clones expressing maltogenic alpha-amylase activity to be identified.

30 Alternatively, the DNA sequence encoding the enzyme may be prepared synthetically by established standard methods, e.g. the phosphoroamidite method described by S.L. Beaucage and M.H. Caruthers (1981) or the method described by Matthes et al. (1984). In the phosphoroamidite method, oligonucleotides are synthesized, e.g. in an automatic DNA synthesizer, purified, annealed, ligated and cloned
35 in appropriate vectors.

Finally, the DNA sequence may be of mixed genomic and synthetic origin, mixed synthetic and cDNA origin or mixed genomic and cDNA origin, prepared by ligating fragments of synthetic, genomic or cDNA origin, wherein the fragments correspond to various parts of the entire DNA sequence, in accordance with techniques
5 well known in the art. The DNA sequence may also be prepared by polymerase chain reaction (PCR) using specific primers, for instance as described in US 4,683,202 or R.K. Saiki et al. (1988).

Site-directed Mutagenesis

Once a maltogenic alpha-amylase-encoding DNA sequence has been isolated,
10 and desirable sites for modification identified, modifications may be introduced using synthetic oligonucleotides. These oligonucleotides contain nucleotide sequences flanking the desired modification sites; mutant nucleotides are inserted during oligo-nucleotide synthesis. In a specific method, a single-stranded gap of DNA, bridging the maltogenic alpha-amylase-encoding sequence, is created in a vector carrying the
15 maltogenic alpha-amylase gene. Then the synthetic nucleotide, bearing the desired modification, is annealed to a homologous portion of the single-stranded DNA. The re-maining gap is then filled in with DNA polymerase I (Klenow fragment) and the construct is ligated using T4 ligase. A specific example of this method is described in Morinaga et al. (1984): US 4,760,025 discloses the introduction of oligonucleotides
20 encoding multiple modifications by performing minor alterations of the cassette. How-ever, an even greater variety of modifications can be introduced at any one time by the Morinaga method because a multitude of oligonucleotides, of various lengths, can be introduced.

Another method of introducing modifications into a maltogenic alpha-amylase-
25 encoding DNA sequences is described in Nelson and Long (1989). It involves a 3-step generation of a PCR fragment containing the desired modification introduced by using a chemically synthesized DNA strand as one of the primers in the PCR reactions. From the PCR-generated fragment, a DNA fragment carrying the modification may be isolated by cleavage with restriction endonucleases and reinserted into an expression
30 plasmid.

Random Mutagenesis

Random mutagenesis is suitably performed either as localised or region-specific random mutagenesis in at least three parts of the gene translating to the amino acid sequence shown in question, or within the whole gene.
35   The random mutagenesis of a DNA sequence encoding a parent maltogenic alpha-amylase may be conveniently performed by use of any method known in the art.

In relation to the above, a further aspect of the present invention relates to a method for generating a variant of a parent Novamyl-like α-amylase, wherein the variant exhibits increased stability at low pH and at low calcium concentration relative to the parent, the method comprising:

5   (a) subjecting a DNA sequence encoding the parent Novamyl-like α-amylase to random mutagenesis,

(b) expressing the mutated DNA sequence obtained in step (a) in a host cell, and

(c) screening for host cells expressing a Novamyl-like α-amylase variant which
10 has an altered property relative to the parent Novamyl-like α-amylase.

Step (a) of the above method of the invention is preferably performed using doped primers, as described in the working examples herein (vide infra).

For instance, the random mutagenesis may be performed by use of a suitable physical or chemical mutagenizing agent, by use of a suitable oligonucleotide, or by
15 subjecting the DNA sequence to PCR generated mutagenesis. Furthermore, the random mutagenesis may be performed by use of any combination of these mutagenizing agents. The mutagenizing agent may, e.g., be one which induces transitions, transversions, inversions, scrambling, deletions, and/or insertions.

Examples of a physical or chemical mutagenizing agent suitable for the present
20 purpose include ultraviolet (UV) irradiation, hydroxylamine, N-methyl-N'-nitro-N-nitrosoguanidine (MNNG), O-methyl hydroxylamine, nitrous acid, ethyl methane sulphonate (EMS), sodium bisulphite, formic acid, and nucleotide analogues. When such agents are used, the mutagenesis is typically performed by incubating the DNA sequence encoding the parent enzyme to be mutagenized in the presence of the
25 mutagenizing agent of choice under suitable conditions for the mutagenesis to take place, and selecting for mutated DNA having the desired properties.

When the mutagenesis is performed by the use of an oligonucleotide, the oligonucleotide may be doped or spiked with the three non-parent nucleotides during the synthesis of the oligonucleotide at the positions which are to be changed. The
30 doping or spiking may be done so that codons for unwanted amino acids are avoided. The doped or spiked oligonucleotide can be incorporated into the DNA encoding the maltogenic alpha-amylase enzyme by any published technique, using e.g. PCR, LCR or any DNA polymerase and ligase as deemed appropriate.

Preferably, the doping is carried out using "constant random doping", in which
35 the percentage of wild-type and modification in each position is predefined. Furthermore, the doping may be directed toward a preference for the introduction of certain nucleotides, and thereby a preference for the introduction of one or more specific amino acid residues. The doping may be made, e.g., so as to allow for the

introduction of 90% wild type and 10% modifications in each position. An additional consideration in the choice of a doping scheme is based on genetic as well as protein-structural constraints. The doping scheme may be made by using the DOPE program which, *inter alia*, ensures that introduction of stop codons is avoided.

5   When PCR-generated mutagenesis is used, either a chemically treated or non-treated gene encoding a parent maltogenic alpha-amylase enzyme is subjected to PCR under conditions that increase the misincorporation of nucleotides (Deshler 1992; Leung et al., Technique, Vol.1, 1989, pp. 11-15).

          A mutator strain of *E. coli* (Fowler et al., Molec. Gen. Genet., 133, 1974, pp.
10   179-191), *S. cereviseae* or any other microbial organism may be used for the random mutagenesis of the DNA encoding the maltogenic alpha-amylase by, e.g., transforming a plasmid containing the parent enzyme into the mutator strain, growing the mutator strain with the plasmid and isolating the mutated plasmid from the mutator strain. The mutated plasmid may be subsequently transformed into the expression organism.

15   The DNA sequence to be mutagenized may be conveniently present in a genomic or cDNA library prepared from an organism expressing the parent maltogenic alpha-amylase. Alternatively, the DNA sequence may be present on a suitable vector such as a plasmid or a bacteriophage, which as such may be incubated with or other-wise exposed to the mutagenising agent. The DNA to be mutagenized may also be
20   present in a host cell either by being integrated in the genome of said cell or by being present on a vector harbored in the cell. Finally, the DNA to be mutagenized may be in isolated form. It will be understood that the DNA sequence to be subjected to random mutagenesis is preferably a cDNA or a genomic DNA sequence.

          In some cases it may be convenient to amplify the mutated DNA sequence
25   prior to performing the expression step b) or the screening step c). Such amplification may be performed in accordance with methods known in the art, the presently preferred method being PCR-generated amplification using oligonucleotide primers prepared on the basis of the DNA or amino acid sequence of the parent enzyme.

          Subsequent to the incubation with or exposure to the mutagenising agent, the
30   mutated DNA is expressed by culturing a suitable host cell carrying the DNA sequence under conditions allowing expression to take place. The host cell used for this purpose may be one which has been transformed with the mutated DNA sequence, optionally present on a vector, or one which was carried the DNA sequence encoding the parent enzyme during the mutagenesis treatment. Examples of suitable host cells are the
35   following: gram positive bacteria such as *Bacillus subtilis*, *Bacillus licheniformis*, *Bacillus lentus*, *Bacillus brevis*, *Bacillus stearothermophilus*, *Bacillus alkalophilus*, *Bacillus amyloliquefaciens*, *Bacillus coagulans*, *Bacillus circulans*, *Bacillus lautus*, *Bacillus*

megaterium, Bacillus thuringiensis, Streptomyces lividans or *Streptomyces murinus*; and gram negative bacteria such as *E. coli.*

The mutated DNA sequence may further comprise a DNA sequence encoding functions permitting expression of the mutated DNA sequence.

5 Localized random mutagenesis

The random mutagenesis may be advantageously localised to a part of the parent maltogenic alpha-amylase in question. This may, e.g., be advantageous when certain regions of the enzyme have been identified to be of particular importance for a given property of the enzyme, and when modified are expected to result in a variant
10 having improved properties. Such regions may normally be identified when the tertiary structure of the parent enzyme has been elucidated and related to the function of the enzyme.

The localized, or region-specific, random mutagenesis is conveniently performed by use of PCR generated mutagenesis techniques as described above or
15 any other suitable technique known in the art. Alternatively, the DNA sequence encoding the part of the DNA sequence to be modified may be isolated, e.g., by insertion into a suitable vector, and said part may be subsequently subjected to mutagenesis by use of any of the mutagenesis methods discussed above.

For region-specific random mutagenesis with a view to improving the stability of
20 calcium binding of a parent maltogenic alpha-amylase, codon positions corresponding to the following amino acid residues from the amino acid sequence set forth in SEQ ID NO: 1 may appropriately be targeted:

Residues:Regions:

16-33, 35-36, 40: 16-40

25   46-54, 56: 46-56

73-81: 73-81

87-89, 91, 93-96, 99-105, 109: 87-109

129-134, (145, 150): 129-134

167-172, 174, 177, 180-189: 167-189

30   196-202, 206-210: 196-210

228-235, 237: 228-237

378

637

With a view to achieving improved binding of a substrate, i.e., improved binding
35 of a carbohydrate species, such as amylose or amylopectin, by a maltogenic alpha-amylase variant with a modified, e.g. higher, substrate specificity and/or a modified, e.g. higher, specificity with respect to cleavage, i.e. hydrolysis, of the substrate, it appears

that the following codon positions in the following regions of the amino acid sequence shown in SEQ ID NO: 1, may particularly appropriately be targeted for modification by region-specific mutagenesis:

    70-97, 127-143, 174-198, 226-233, 255-270, 282-292, 324-331, 370-376.

5   For region-specific random mutagenesis with a view to altering the substrate specificity and/or the pH dependent activity profile, the following regions of SEQ ID NO: 1 may be targeted: 70-97, 174-198.

    The following regions may be targeted with a view to improving the thermostability: 70-109, 167-200.

10 <u>General method for random mutagenesis by use of the DOPE program</u>

    The random mutagenesis may be carried out by the following steps:

    1.Select regions of interest for modification in the parent enzyme

    2.Decide on mutation sites and nonmutated sites in the selected region

    3.Decide on which kind of mutations should be carried out, e.g. with respect to
15 the desired stability and/or performance of the variant to be constructed

    4.Select structurally reasonable mutations

    5.Adjust the residues selected by step 3 with regard to step 4.

    6.Analyse by use of a suitable dope algorithm the nucleotide distribution.

    7.If necessary, adjust the wanted residues to genetic code realism, e.g. taking
20 into account constraints resulting from the genetic code, e.g. in order to avoid introduction of stop codons; the skilled person will be aware that some codon combinations cannot be used in practice and will need to be adapted

    8.Make primers

    9.Perform random mutagenesis by use of the primers

25   10.Select resulting α-amylase variants by screening for the desired improved properties.

    Suitable dope algorithms for use in step 6 are well known in the art. One such algorithm is described by Tomandl, D. et al., 1997, Journal of Computer-Aided Molecular Design 11:29-38. Another algorithm is DOPE (Jensen, LJ, Andersen, KV,
30 Svendsen, A, and Kretzschmar, T (1998) Nucleic Acids Research 26:697-702).

Expression of maltogenic alpha-amylase variants

    The construction of the variant of interest is accomplished by cultivating a microorganism comprising a DNA sequence encoding the variant under conditions
35 which are conducive for producing the variant, and optionally subsequently recovering the variant from the resulting culture broth. This is described in detail further below.

According to the invention, a DNA sequence encoding the variant produced by methods described above, or by any alternative methods known in the art, can be expressed, in the form of a protein or polypeptide, using an expression vector which typically includes control sequences encoding a promoter, operator, ribosome binding
5 site, translation initiation signal, and, optionally, a repressor gene or various activator genes.

The recombinant expression vector carrying the DNA sequence encoding an maltogenic alpha-amylase variant of the invention may be any vector which may conveniently be subjected to recombinant DNA procedures, and the choice of vector
10 will often depend on the host cell into which it is to be introduced. Thus, the vector may be an autonomously replicating vector, i.e. a vector which exists as an extrachromo-somal entity, the replication of which is independent of chromosomal replication, e.g. a plasmid, a bacteriophage or an extrachromosomal element, minichromosome or an artificial chromosome. Alternatively, the vector may be one which, when introduced into
15 a host cell, is integrated into the host cell genome and replicated together with the chromosome(s) into which it has been integrated.

In the vector, the DNA sequence should be operably connected to a suitable promoter sequence. The promoter may be any DNA sequence which shows transcrip-tional activity in the host cell of choice and may be derived from genes encoding pro-
20 teins either homologous or heterologous to the host cell. Examples of suitable pro-moters for directing the transcription of the DNA sequence encoding a maltogenic alpha-amylase variant of the invention, especially in a bacterial host, are the promoter of the *lac* operon of *E.coli*, the *Streptomyces coelicolor* agarase gene *dagA* promoters, the promoters of the *Bacillus licheniformis* α-amylase gene (*amyL*), the promoters of
25 the *Bacillus stearothermophilus* maltogenic amylase gene (*amyM*), the promoters of the *Bacillus amyloliquefaciens* α-amylase (*amyQ*), the promoters of the *Bacillus subtilis* xylA and xylB genes, etc. For transcription in a fungal host, examples of useful promo-ters are those derived from the gene encoding *A. oryzae* TAKA amylase, *Rhizomucor miehei* aspartic proteinase, *A. niger* neutral α-amylase, *A. niger* acid stable α-amylase,
30 *A. niger* glucoamylase, *Rhizomucor miehei* lipase, *A. oryzae* alkaline protease, *A. oryzae* triose phosphate isomerase or *A. nidulans* acetamidase.

The expression vector of the invention may also comprise a suitable transcription terminator and, in eukaryotes, polyadenylation sequences operably connected to the DNA sequence encoding the maltogenic alpha-amylase variant of the
35 invention. Termination and polyadenylation sequences may suitably be derived from the same sources as the promoter.

The vector may further comprise a DNA sequence enabling the vector to replicate in the host cell in question. Examples of such sequences are the origins of replication of plasmids pUC19, pACYC177, pUB110, pE194, pAMB1 and pIJ702.

The vector may also comprise a selectable marker, e.g. a gene the product of
5 which complements a defect in the host cell, such as the *dal* genes from *B. subtilis* or *B. licheniformis*, or one which confers antibiotic resistance such as ampicillin, kanamycin, chloramphenicol or tetracycline resistance. Furthermore, the vector may comprise *Aspergillus* selection markers such as amdS, argB, niaD and sC, a marker giving rise to hygromycin resistance, or the selection may be accomplished by co-
10 transformation, e.g. as described in WO 91/17243.

While intracellular expression may be advantageous in some respects, e.g. when using certain bacteria as host cells, it is generally preferred that the expression is extracellular. In general, the *Bacillus* α-amylases mentioned herein comprise a preregion permitting secretion of the expressed protease into the culture medium. If desir-
15 able, this preregion may be replaced by a different preregion or signal sequence, conveniently accomplished by substitution of the DNA sequences encoding the respective preregions.

The procedures used to ligate the DNA construct of the invention encoding maltogenic alpha-amylase variant, the promoter, terminator and other elements,
20 respectively, and to insert them into suitable vectors containing the information necessary for replication, are well known to persons skilled in the art (cf., for instance, Sambrook et al. (1989)).

The cell of the invention, either comprising a DNA construct or an expression vector of the invention as defined above, is advantageously used as a host cell in the
25 recombinant production of a maltogenic alpha-amylase variant of the invention. The cell may be transformed with the DNA construct of the invention encoding the variant, conveniently by integrating the DNA construct (in one or more copies) in the host chromosome. This integration is generally considered to be an advantage as the DNA sequence is more likely to be stably maintained in the cell. Integration of the DNA con-
30 structs into the host chromosome may be performed according to conventional methods, e.g. by homologous or heterologous recombination. Alternatively, the cell may be transformed with an expression vector as described above in connection with the different types of host cells.

The cell of the invention may be a cell of a higher organism such as a mammal
35 or an insect, but is preferably a microbial cell, e.g. a bacterial or a fungal (including yeast) cell.

Examples of suitable bacteria are gram positive bacteria such as *Bacillus subtilis, Bacillus licheniformis, Bacillus lentus, Bacillus brevis, Bacillus stearothermo-*

philus, *Bacillus alkalophilus*, *Bacillus amyloliquefaciens*, *Bacillus coagulans*, *Bacillus circulans*, *Bacillus lautus*, *Bacillus megaterium*, *Bacillus thuringiensis*, or *Streptomyces lividans* or *Streptomyces murinus*, or gram negative bacteria such as *E.coli*. The trans- formation of the bacteria may, for instance, be effected by protoplast transformation or
5 by using competent cells in a manner known *per se*.

The yeast organism may favourably be selected from a species of *Saccharomyces* or *Schizosaccharomyces*, e.g. *Saccharomyces cerevisiae*. The filamentous fungus may advantageously belong to a species of *Aspergillus*, e.g. *Aspergillus oryzae* or *Aspergillus niger*. Fungal cells may be transformed by a process
10 involving protoplast formation and transformation of the protoplasts followed by regen- eration of the cell wall in a manner known *per se*. A suitable procedure for transformation of *Aspergillus* host cells is described in EP 238 023.

In a yet further aspect, the present invention relates to a method of producing a maltogenic alpha-amylase variant of the invention, which method comprises cultivating
15 a host cell as described above under conditions conducive to the production of the variant and recovering the variant from the cells and/or culture medium.

The medium used to cultivate the cells may be any conventional medium suitable for growing the host cell in question and obtaining expression of the maltogenic alpha-amylase variant of the invention. Suitable media are available from commercial
20 suppliers or may be prepared according to published recipes (e.g. as described in cata- logues of the American Type Culture Collection).

The maltogenic alpha-amylase variant secreted from the host cells may con- veniently be recovered from the culture medium by well-known procedures, including separating the cells from the medium by centrifugation or filtration, and precipitating
25 proteinaceous components of the medium by means of a salt such as ammonium sulfate, followed by the use of chromatographic procedures such as ion exchange chromatography, affinity chromatography, or the like.

Testing of maltogenic alpha-amylase variants

Maltogenic alpha-amylase variants produced by any of the methods described
30 above may be tested, either prior to or after purification, for amylolytic activity in a screening assay which measures the ability of the variant to degrade starch. The screening in step 10 in the above-mentioned random mutagenesis method of the invention may be conveniently performed by use of a filter assay based on the following procedure: A microorganism capable of expressing the mutated maltogenic alpha-
35 amylase of interest is incubated on a suitable medium and under suitable conditions for secretion of the enzyme, the medium being covered with two filters comprising a protein-binding filter placed under a second filter exhibiting a low protein binding

capability. The microorganism is grown on the second, top filter. Subsequent to the incubation, the bottom protein-binding filter comprising enzymes secreted from the microorganism is separated from the second filter comprising the microorganism. The protein-binding filter is then subjected to screening for the desired enzymatic activity,

5 and the corresponding microbial colonies present on the second filter are identified. The first filter used for binding the enzymatic activity may be any protein-binding filter, e.g., nylon or nitrocellulose. The second filter carrying the colonies of the expression organism may be any filter that has no or low affinity for binding proteins, e.g., cellulose acetate or Durapore™.

10 Screening consists of treating the first filter to which the secreted protein is bound with a substrate that allows detection of the α-amylase activity. The enzymatic activity may be detected by a dye, fluorescence, precipitation, pH indicator, IR-absorbance or any other known technique for detection of enzymatic activity. The detecting compound may be immobilized by any immobilizing agent e.g. agarose, agar,

15 gelatine, polyacrylamide, starch, filter paper, cloth; or any combination of immobilizing agents. For example, α-amylase activity can be detected by Cibacron Red labelled amylopectin, which is immobilized in agarose. α-amylase activity on this substrate produces zones on the plate with reduced red color intensity.

To screen for variants with increased stability, the filter with bound maltogenic

20 alpha-amylase variants can be pretreated prior to the detection step described above to inactivate variants that do not have improved stability relative to the parent maltogenic alpha-amylase. This inactivation step may consist of, but is not limited to, incubation at elevated temperatures in the presence of a buffered solution at any pH from pH 2 to 12, and/or in a buffer containing another compound known or thought to contribute to

25 altered stability e.g., surfactants, EDTA, EGTA, wheat flour components, or any other relevant additives. Filters so treated for a specified time are then rinsed briefly in deionized water and placed on plates for activity detection as described above. The conditions are chosen such that stabilized variants show increased enzymatic activity relative to the parent after incubation on the detection media.

30 To screen for variants with altered thermostability, filters with bound variants are incubated in buffer at a given pH (e.g., in the range from pH 2-12) at an elevated temperature (e.g., in the range from 50°-110°C) for a time period (e.g., from 1-20 minutes) to inactivate nearly all of the parent maltogenic alpha-amylase, rinsed in water, then placed directly on a detection plate containing immobilized Cibacron Red

35 labelled amylopectin and incubated until activity is detectable. Similarly, pH dependent stability can be screened for by adjusting the pH of the buffer in the above inactivation step such that the parent maltogenic alpha-amylase is inactivated, thereby allowing detection of only those variants with increased stability at the pH in question. To screen

for variants with increased calcium-dependent stability calcium chelators, such as ethylene glycol-bis(ß-aminoethyl ether) N,N,N',N'-tetraacetic acid (EGTA), is added to the inactivation buffer at a concentration such that the parent maltogenic alpha-amylase is inactivated under conditions further defined, such as buffer pH, temperature or a
5 specified length of incubation.

The variants of the invention may be suitably tested by assaying the starch-degrading activity of the variant, for instance by growing host cells transformed with a DNA sequence encoding a variant on a starch-containing agarose plate and identifying starch-degrading host cells as described above. Further testing in regard to altered
10 properties, including specific activity, substrate specificity, cleavage pattern, thermoactivation, thermostability, pH dependent activity or optimum, pH dependent stability, temperature dependent activity or optimum, transglycosylation activity, stability, and any other parameter of interest, may be performed on purified variants in accordance with methods known in the art as described below.

15 Degradation of β-limit dextrin by maltogenic alpha-amylase:

Another important parameter in the evaluation of the substrate specificity of maltogenic alpha-amylase variants may be the degree to which such enzymes are capable of degrading starch that has been exhaustively treated with the exoglycosylase β-amylase. To screen for variants which show patterns of degradation on such a
20 substrate differing from the patterns produced by the parent maltogenic alpha-amylase the following assay is performed: β-limit dextrin is prepared by incubating 25 ml 1% amylopectin in McIlvane buffer (48.5 mM citrate and 193 mM sodium phosphate pH 5.0) with 24 µg/ml β-amylase overnight at 30°C. Unhydrolysed amylopectin (i.e., β-limit dextrin) is precipitated with 1 volume 98% ethanol, washed and redissolved in water. 1
25 ml β-limit dextrin is incubated with 18 µl enzymes (at 2.2 mg/ml) and 100 µl 0.2 M citrate-phosphate pH 5.0 for 2 hrs at 30°C and analysed by HPLC as described above. Total hydrolysis of β-limit dextrin is carried out in 2M HCl at 95°C. The concentration of reducing ends is measured by methods known in the art.

Calcium binding affinity

30   Unfolding of maltogenic alpha-amylases by exposure to heat or to denaturants such as guanidine hydrochloride is accompanied by a decrease in fluorescence, and oss of calcium ions leads to unfolding. Thus, the affinity of a maltogenic alpha-amylase variant for calcium can be measured by fluorescence measurements before and after incubation of the variant (e.g., at a concentration of 10 mg/ml) in a buffer (e.g., 50 mM
35 HEPES, pH 7) with different concentrations of calcium (e.g., in the range from

1 mM-100 mM) or of EGTA (e.g., in the range from 1-1000 mM) for a sufficiently long period of time (such as 22 hours at 55°C).

The measured fluorescence, F, is composed of contributions form the folded and unfolded forms of the enzyme. The following equation can be derived to describe the dependence of F on calcium concentration ([Ca]):

$$F = [Ca]/(K_{diss} + [Ca])(a_N - b_N log([Ca])) +$$

$$K_{diss}/(K_{diss} + [Ca])(a_U - b_U log([Ca]))$$

where a_N is the fluorescence of the native (folded) form of the enzyme, b_N is the linear dependence of a_N on the logarithm of the calcium concentration (as observed experimentally), a_U is the fluorescence of the unfolded form and b_U is the linear dependence of a_U on the logarithm of the calcium concentration. K_{diss} is the apparent calcium binding constant for an equilibrium process as follows:

$$K_{diss}$$

N-Ca« U + Ca(N = native enzyme; U = unfolded enzyme)

In fact, unfolding proceeds extremely slowly and is irreversible. The rate of unfolding is dependent on calcium concentration, and such dependency for a given enzyme provides a measure of the calcium binding affinity of the enzyme. By defining a standard set of reaction conditions (e.g., 22 hours at 55°C), a meaningful comparison of K_{diss} for different maltogenic alpha-amylase variants can be made.

Industrial Applications

The maltogenic alpha-amylase variants of the invention possesses valuable properties which may be advantageously used in various industrial applications. In particular, the enzyme finds potential application for retarding or preventing retrogradation, and thus the staling, of starch based food such as common in the baking industry.

The variant may be used for the preparation of bread and other bread products in accordance with conventional techniques known in the art.

It is believed that the modification of the starch fraction by use of the present invention results in increased volume in baked products and improved organoleptic qualities, such as flavour, mouth feel, palatability, aroma and crust colour.

The maltogenic alpha-amylase variant may be used as the only enzyme or as a major enzymatic activity in combination with one or more additional enzymes, such as xylanase, lipase, glucose oxidase and other oxidoreductases, or an amylolytic enzyme.

The enzyme variants of the invention also find industrial applicability as a component in washing, dishwashing and hard-surface cleaning detergent compositions. Some variants are particularly useful in a process for the manufacture of

linear oligosaccharides, or in the production of sweeteners and ethanol from starch, and/or for textile desizing. Conditions for conventional starch conversion processes, including starch liquefaction and/or saccharification processes, are described in, e.g., US 3,912,590 and in EP patent publications Nos. 252,730 and 63,909.

5   The invention is further illustrated with reference to the following examples which are not intended to be in any way limiting to the scope of the invention as claimed.

Determination of maltogenic amylase in MANU

One Maltogenic Amylase Novo Unit (MANU) is the amount of enzyme which 10 under standard will cleave one μmol maltotriose per minute. The standard conditions are 10 mg/ml maltotriose, 37°C, pH 5.0, 30 minutes reaction time.

The pH dependence is found by repeating this measurement at the same conditions, but at different pH values.

EXAMPLES

15 **Example 1: Construction of a variant of Novamyl with altered pH dependent activity**

Novamyl is expressed in *Bacillus subtilis* from a plasmid denoted herein as pLBei010. This plasmid contains *amyM* in which the expression of *amyM* is directed by its own promoter and the complete gene encoding Novamyl, e.g., as contained in the 20 strain DSM 11837. The plasmid contains the origin of replication, *ori*, from plasmid pUB110 and an kanamycin resistance marker for selection purposes. pLBei010 is shown in Fig. 1.

Primer sequences

Site directed mutants of Novamyl were constructed by the megaprimer method 25 essentially as described by Kammann et al. (1989). Briefly, a mutagenic oligonucleotide primer is used together in a PCR reaction with a suitable opposite DNA strand end primer to create a preliminary PCR product. This product is then used as a megaprimer together with another opposite DNA strand end primer to create a double-stranded DNA product. The product of the final PCR reaction was routinely used to replace a 30 corresponding DNA fragment in the pLBei010 plasmid by standard cloning procedures. Mutants were transformed directly into *Bacillus subtilis* strain SHa273, a derivative of *Bacillus subtilis* 168 which is apr⁻, npr⁻, amyE⁻, amyR2⁻ and prepared by methods known in the art.

Oligonucleotide primers used in the construction of described variants are as listed below:

Variant Sequence (5'→ 3')

F188H: SEQ ID NO: 3

5   F188E: SEQ ID NO: 4

F284E: SEQ ID NO: 5

F284D: SEQ ID NO: 6

F284K: SEQ ID NO: 7

N327D: SEQ ID NO: 8

10   Variant Sequence (3'→ 5')

T288K: SEQ ID NO: 9

T288R: SEQ ID NO: 10

Aspartate variants of F284, T288 and N327 were obtained using primer A189 (SEQ ID NO: 11) and B649 (SEQ ID NO: 12) as end-primers.

15   F188-variants F188L, T189Y were obtained using primer A82 (SEQ ID NO: 13) and B346 (SEQ ID NO: 14) as end-primers.

PCR products with the desired modification(s) were purified, digested with appropriate enzymes, separated by agarose gel electrophoresis and extracted, ethanol precipitated in the presence of glycogen, resuspended in H_2O, ligated to pLBei010

20 which had been digested with the same appropriate enzymes, and transformed into *Bacillus subtilis* SHa273. Transformants were checked for size by colony PCR and for the insertion or removal of specific restriction sites by restriction enzyme digestion. Positive colonies were verified by DNA sequencing methods as described in the art.

Fermentation

25   The *B. subtilis* SHa273 mutant clones were grown overnight on LB-Kana (10 µg/ml)-Starch plates at 37°C. The colonies from the plate were resuspended in 10 ml Luria broth. One-sixth of each of the suspensions were inoculated into a 500 ml shake flasks containing 100 ml PS-1 media, a soy meal/sucrose-based media, kanamycin for a final concentration of 10 µg/ml and 100 µl 5M NaOH. The pH was adjusted to 7.5 with

30 NaOH before inoculation. The cultures were incubated for five days at 30°C with shaking at 270-300 rpm.

Enzyme Purification

Large particles from the media were removed by flocculation before affinity chromatography. Superfloc C521 (American Cyanmide Company) was used as the

35 cationic flocculant and Superfloc A130 (American Cyanmide Company) as the anionic flocculant.

The culture suspension was diluted 1:1 with deionized water and the pH was adjusted to approx. 7.5. A volume of 0.01 ml of $50^w/_w\%$ $CaCl_2$ per ml diluted culture was added during stirring. A volume of 0.015 ml of $20^w/_w\%$ Na-aluminate per ml diluted culture was titrated with 20% formic acid, while keeping the pH between 7 and 8. While
5   stirring 0.025 ml $10^v/_v\%$ of C521 per ml diluted culture was added, followed by 0.05 ml $1^w/_v\%$ A130 per ml diluted culture, or until flocculation was observed. The solution was centrifuged at 4500 rpm for 30 minutes. Filtration was performed using a filter of pore size of 0.45 μm to exclude larger particles and any remaining bacteria. The filtered solution was stored at -20°C.

10   ## Immobilization of α-cyclodextrin to DSV-agarose

One hundred mg of α-cylcodextrin of molecular weight 972.86g/mol (Fluka 28705) was dissolved in 20 ml coupling buffer (0.5M Na_2CO_3, pH 11). Ten ml of DSV-agarose (Mini-Leak, Medium 10-20 mmol/l of divinyl sulfone activated agarose (Kem-En-Tec) was washed thoroughly with deionized water, then dried by suction and
15   transferred to the a-cyclodextrin solution. After the mixture had stirred for 24 hr at ambient temperature, the gel was washed with deionized water, followed by 0.5M $KHCO_3$. The gel was transferred to the blocking buffer (20ml 0.5M $KHCO_3$ + 1ml mercaptoethanol), stirred for 2 hr at ambient temperature, then washed with deionized water.

20   ## Affinity chromatography

The variants were purified by affinity chromatography using the Pharmacia FPLC System. A 0.04 volume of 1M Na-acetate pH 5 was added to the filtrate obtained by flocculation to adjust pH and $CaCl_2$ was added to a final concentration of 10^{-10} M. The solution was filtered and degassed. A Pharmacia XK16 column was prepared with
25   ten ml of the immobilised α-cyclodextrin, then equilibrated in the equilibration buffer (25 mM Na-acetate pH 5) by washing with approximately 10 times the column volume. The filtrate was applied to the XK16 column, which was then washed with the equilibration buffer until protein could no longer be detected in the washing buffer. The column was washed with the equilibration buffer containing 0.5M NaCl to elute nonspecific material,
30   followed by another wash with 2-3 times the column volume of the equilibration buffer. All washings were performed using a flow rate of 10ml/min. Specifically bound material was eluted using a solution of 2% α-cyclodextrin in the wash buffer and collected using the Pharmacia Liquid Chromatography Collector LCC-500 Plus using a flow rate of 5 ml/min.

Example 2: pH dependent activity of variants

The variants prepared in the preceding Example were tested for activity at various pH values as follows.

A colorimetric glucose oxidase-peroxidase assay for liberated glucose from
5 maltotriose or amylopectin was used to determine the pH activity profiles of the enzyme variants (Glucose/GOD-Perid® Method, Boehringer Mannheim, Indianapolis IN). Activity was assayed in a buffer of 25 mM citrate-phosphate, 0.1mM $CaCl_2$ at pH values of 2, 2.5, 3, 3.5, 4, 4.5, 5, 5.5, 6, 6.5, 7, 7.5, 8 and 8.6. The buffer pH was adjusted using NaOH and enzymes were diluted in 25 mM citrate-phosphate buffer pH 5.
10 Measurements were taken in duplicate to obtain an average value. All values are relative to the pH at which the highest level of activity is seen.

The results, shown in the table below, indicate that each of the variants has an alteration in the pH dependent activity profile when compared to the parent Novamyl®. The highest level of activity for each variant is designated 100% and the activity of that
15 variant measured at the other indicated pH values is a relative percentage of that maximum.

Modifications	pH													
	2.0	2.5	3.0	3.5	4.0	4.5	5.0	5.5	6.0	6.5	7.0	7.5	8.0	8.6
None (parent)	0	0	0	8	47	80	100	95	91	80	66	39	35	30
F188H	1	0	0	1	3	29	77	99	100	88	59	39	31	27
F188E	0	0	0	2	27	62	89	100	93	71	46	28	20	18
T288R	0	0	0	8	51	77	94	100	86	73	50	34	27	12
N327D	1	1	7	27	67	95	100	98	77	33	19	11	5	0

Further, a number of Novamyl variants were tested for activity at pH 4.0 and
20 5.0, taking the activity of Novamyl at the same pH as 100 %. The activity was determined by hydrolysis of maltotriose (10 mg/ml) at 60°C, 50 mM sodium acetate, 1 mM $CaCl_2$. The results are expressed as the ratio between activity at pH 5.0 and pH 4.0:

Modifications	pH 5.0/pH 4.0
N131D	0.24
I174Q	0.31

G397P	0.40
H103Y	0.40
Δ 262-266	0.47
S32Q	0.53
S32D	0.55
T142A+ D261G	0.62
G370N+ N371G	0.66
S32N	0.68
N176S	0.79
D17E	0.80
None (parent)	1
Δ 191	1.39
192-A-193	1.61
I174E	1.80
192-A-G-193	1.90
Δ 192	2.22

The results demonstrate that variants with a higher or lower pH optimum can be obtained according to the invention.

Example 3: Thermostability of variants

Incubation at 80°C

The thermostability of a number of Novamyl variants was tested by incubating an aqueous solution at 80°C and pH 4.3 and measuring the residual amylase activity at various times. The parent enzyme, Novamyl, was included for comparison. The results are expressed as residual activity at various times in percent of initial activity:

Modifications	**0**	**5 min.**	**10 min.**	**15 min.**	**20 min.**	**25 min.**
None (parent)	100	23	9	3	1	0
F188L+ V336L+ T525A	100	63	49	48	52	47
F188I+ Y422F+ I660V	100	71	60	51	43	38
N115D+ F188L	100	73	60	51	44	39

A30D+ K40R+ D261G	100	38	24	15	13	10
T142A+ N327S+ K425E+ K520R+ N595I	100	47	39	25	19	11
F188L+ D261G+ T288P	100	60	67	66	63	67
K40R+ F188L+ D261G+ A483T	100	56	48	40	36	30
T288K	100	64	31	18	7	4

The above data show a clearly improved thermostability for the variants compared to the parent amylase.

<u>Incubation at 85°C with calcium</u>

The Novamyl variant S32E was tested by incubation with 1 mM Ca^{++} at 85°C for 15 minutes. The variant showed a residual activity of 48 % whereas the parent enzyme (Novamyl) showed 32 % residual activity at the same conditions.

<u>DSC</u>

Further, the thermostability was tested for some Novamyl variants by DSC (differential scanning calorimetry) at pH 4.3 or 5.5. Again, the parent amylase was included for comparison. The results are expressed as the denaturation temperature (Tm):

Modifications	**Tm at pH 4.3**	**Tm at pH 5.5**
None (parent)	79°C	88°C
N115D+ F188L	86°C	92°C
T142A+ N327S+ K425E+ K520R+ N595I	not determined	93°C

The results show improved thermostability for both variants.

Example 4: Specific activity of variants

Amylase activity was determined by a colorimetric measurement after action on Phadebas tablets at pH 5.0 and 60°C. The resutls for two Novamyl variants, relative to Novamyl were as follows:

Modifications	Relative amylase activity
None (parent)	100
192-A-193	110
Δ (191-195)	300

The specific activity was further tested by action on maltotriose at pH 4.0, 60°C by the MANU method described above. The results showed that the variant G370N, N371G has a maltotriose activity of 106 % compared to Novamyl.

5 **Example 5: Inhibition of retrogradation**

The efficiency of Novamyl and Novamyl variants to inhibit retrogradation was determined as follows:

730 mg of 50 % (w/w) amylopectin slurry in 0.1 M sodium acetate, at a selected pH (3.7, 4.3 or 5.5) was mixed with 20 μl of an enzyme sample, and the mixture was
10 incubated in a sealed ampoule for 1 hour at 40°C, followed by incubation at 100 °C for 1 hour in order to gelatinize the samples. The sample was then aged for 7 days at room temperature to allow recrystallization of the amylopectin. A control without enzyme was included.

After aging, DSC was performed on the sample by scanning from 5°C to 95°C
15 at a constant scan rate of 90°C/hour. The area under the first endothermic peak in the thermogram was taken to represent the amount of retrograded amylopectin, and the relative inhibition of retrogradation was taken as the area reduction (in %) relative to the control without enzyme.

In the table below, the efficiency of the enzyme is expressed as the ratio of the
20 relative inhibition of retrogradation to the enzyme dosage (in MANU/ml):

pH	Modifications	MANU/ml	Relative inhibition	Efficiency
3.7	A30D+ K40R+ D261G	0.23	0.38	1.7
3.7	T142A+ N327S+ K425E+ K520R+ N595I	0.07	0.29	4.1
3.7	None (parent)	0.27	0.38	1.4
4.3	N115D+ F188L	0.01	0.18	18
4.3	None (parent)	0.27	0.43	1.6

5.5	Δ (191-195)+ F188L+ T189Y	0.02	0.12	6
5.5	Δ (191-195)	0.02	0.14	7
5.5	Δ (191-195)	0.05	0.31	6.2
5.5	N115D+ F188L	0.01	0.39	39
5.5	T142A+ D261G	0.14	0.53	3.8
5.5	None (parent)	0.27	0.49	1.8

The results demonstrate that a number of variants are more efficient than the parent amylase to inhibit retrogradation.

Example 6: Anti-staling effect of variants

5   Bread was made by the European Straight Dough method or from sour dough with or without addition of enzymes, and loaves were baked in lidded pans, to avoid volume effects. The bread was allowed to cool for 2 hours, and the texture was analyzed. The remaining loaves were then wrapped in plastic bags and stored at room temperature for texture analysis after 1, 4 and 7 days.

10   The texture analysis of each loaf was done by cutting 4 slices; the force was measured at 25 % compression (P1), at 40% compression (P2) and after keeping 40% compression constant for 30 sec. (P3). P1 was taken as the firmness, and the ratio (P3/P2) was taken as the elasticity of the crumb. The extent of retrogradation after 7 days storage was determined by DSC as described in Example 7.

15 European Straight Dough (pH5.5-6.0)

A Novamyl variant (T142A+ N327S+ K425E+ K520R+ N595I) was tested at dosages in the range of 0-2 mg enzyme/kg flour, and the parent enzyme (Novamyl) was used for comparison.

The results showed that at equal dosages, the variant gives better elasticity
20 (P3/P2) than the parent enzyme after two hours and 1 day. The results after 7 days showed that the variant at dosages of 1-2 mg/kg gave significantly softer crumb (lower firmness, P1) than the parent enzyme at the same dosage. Thus, the variant has a better anti-staling effect throughout a 7-day storage period.

Sour dough (pH approx. 4.5)

25   A Novamyl variant (F188L+ D261G+ T288P) was tested in sour dough, and the parent enzyme (Novamyl) was used for comparison. The following results were

obtained for firmness (P1) after 7 days, elasticity (P3/P2) after 4 and 7 days and retrogradation after 7 days:

Enzyme	Dosage mg/kg flour	Firmness (P1) after 7 days
None	0	2590
Parent	1	2031
	3	1912
	13	1570
Variant	1	1436
	3	1226

Enzyme	Dosage mg/kg flour	Elasticity 4 days	Elasticity 7 days
None	0	0.49	0.47
Parent	1	0.51	0.52
	3	0.53	0.51
	13	0.53	0.51
Variant	1	0.59	0.57
	3	0.57	0.58

Enzyme	Dosage mg/kg flour	Retrogradation, 7 days (relative to control)
None	0	100 %
Parent	1	100 %
	3	63 %
	13	32 %
Variant	1	46 %
	3	20 %

5

The results show that the variant has a markedly improved effect on texture evaluated as firmness and elasticity in sour dough at pH 4.5. A dosage of 1-3 mg/kg of the variant is superior to 13 mg/kg of the parent enzyme on all parameters tested, and the elasticity achieved with the variant cannot be matched by the parent enzyme 5 at any dosage.

Example 7: Cleavage pattern of variants

The cleavage pattern in starch hydrolysis was compared for two variants and the parent enzyme, Novamyl.

The results below indicate % by weight of each oligosacccharide (G1-G8) 10 formed after 24 hours incubation in 1% (w/v) starch using 50 mM sodium acetate, 1 mM $CaCl_2$, pH 5.0 at 50 °C. The oligosaccharides were identified and quantitated using HPLC.

Oligosaccharide	Parent	Δ (191-195)	N115D+ F188L
G8	-	1.7	-
G7	-	2.6	-
G6	-	7.5	1.4
G5	-	10.1	2.1
G4	-	21.1	11.3
G3	-	28.7	10.7
G2	96.5	28.3	61.9
G1	3.5	-	12.6

The results demonstrate a significantly altered cleavage pattern. Novamyl after 15 24 hours produces mainly maltose and virtually no higher oligosaccharides. In contrast, the two variants produce significant amounts of maltotriose and higher oligosaccharides.

REFERENCES CITED

Klein, C., et al., Biochemistry 1992, 31, 8740-8746,
20   Mizuno, H., et al., J. Mol. Biol. (1993) 234, 1282-1283,
Chang, C., et al, J. Mol. Biol. (1993) 229, 235-238,
Larson, S.B., J. Mol. Biol. (1994) 235, 1560-1584,
Lawson, C.L., J. Mol. Biol. (1994) 236, 590-600,
Qian, M., et al., J. Mol. Biol. (1993) 231, 785-799,

Brady, R.L., et al., *Acta Crystallogr. sect. B*, **47**, 527-535,

Swift, H.J., et al., *Acta Crystallogr. sect. B*, **47**, 535-544

A. Kadziola, Ph.D. Thesis: "An alpha-amylase from Barley and its Complex with a Substrate Analogue Inhibitor Studied by X-ray Crystallography", Department of Chemistry University of Copenhagen 1993

MacGregor, E.A., Food Hydrocolloids, 1987, Vol.1, No. 5-6, p.

B. Diderichsen and L. Christiansen, Cloning of a maltogenic α-amylase from *Bacillus stearothermophilus*, FEMS Microbiol. letters: 56: pp. 53-60 (1988)

Hudson et al., Practical Immunology, Third edition (1989), Blackwell Scientific Publications,

Sambrook et al., Molecular Cloning: A Laboratory Manual, 2nd Ed., Cold Spring Harbor, 1989

S.L. Beaucage and M.H. Caruthers, Tetrahedron Letters 22, 1981, pp. 1859-1869

Matthes et al., The EMBO J. 3, 1984, pp. 801-805.

R.K. Saiki et al., Science 239, 1988, pp. 487-491.

Morinaga et al., (1984, Biotechnology 2:646-639)

Nelson and Long, Analytical Biochemistry 180, 1989, pp. 147-151

Hunkapiller et al., 1984, Nature 310:105-111

R. Higuchi, B. Krummel, and R.K. Saiki (1988). A general method of *in vitro* preparation and specific mutagenesis of DNA fragments: study of protein and DNA interactions. *Nucl. Acids Res.* **16**:7351-7367.

Dubnau et al., 1971, J. Mol. Biol. 56, pp. 209-221.

Gryczan et al., 1978, J. Bacteriol. 134, pp. 318-329.

S.D. Erlich, 1977, Proc. Natl. Acad. Sci. 74, pp. 1680-1682.

Boel et al., 1990, Biochemistry 29, pp. 6244-6249.

Kammann, M Laufs, J Schell, J and Gronnenborn, B (1989) Nucleic Acids Research 20:4937-4938.

#	Atom	Res	Chain	ResNo	X	Y	Z	Occ	B	El
1	N	SER	A	1	10.254	56.595	38.175	1.00	15.64	7
2	CA	SER	A	1	11.216	55.462	37.898	1.00	15.87	6
3	C	SER	A	1	12.466	55.723	38.726	1.00	14.53	6
4	O	SER	A	1	12.585	56.773	39.369	1.00	15.99	8
5	CB	SER	A	1	11.527	55.345	36.397	1.00	21.54	6
6	OG	SER	A	1	12.305	56.503	36.045	1.00	20.33	8
7	N	SER	A	2	13.466	54.795	38.551	1.00	18.07	7
8	CA	SER	A	2	14.705	55.061	39.291	1.00	19.33	6
9	C	SER	A	2	15.621	56.069	38.559	1.00	15.87	6
10	O	SER	A	2	16.573	56.563	39.209	1.00	16.73	8
11	CB	SER	A	2	15.490	53.735	39.422	1.00	26.53	6
12	OG	SER	A	2	15.918	53.392	38.123	1.00	21.07	8
13	N	SER	A	3	15.136	56.545	37.384	1.00	12.71	7
14	CA	SER	A	3	15.956	57.522	36.680	1.00	13.38	6
15	C	SER	A	3	15.873	58.916	37.316	1.00	12.57	6
16	O	SER	A	3	16.759	59.749	37.029	1.00	15.22	8
17	CB	SER	A	3	15.434	57.682	35.219	1.00	16.30	6
18	OG	SER	A	3	15.593	56.381	34.568	1.00	23.61	8
19	N	ALA	A	4	14.811	59.222	38.050	1.00	10.88	7
20	CA	ALA	A	4	14.574	60.623	38.384	1.00	11.38	6
21	C	ALA	A	4	15.599	61.115	39.409	1.00	12.81	6
22	CB	ALA	A	4	13.132	60.682	38.956	1.00	12.58	6
23	O	ALA	A	4	15.968	60.306	40.380	1.00	14.28	8
24	N	SER	A	5	16.905	60.780	41.427	1.00	13.21	7
25	CA	SER	A	5	18.163	59.941	41.357	1.00	14.29	6
26	C	SER	A	5	18.053	58.724	41.237	1.00	16.01	6
27	O	SER	A	5	16.218	60.613	42.785	1.00	16.41	8
28	CB	SER	A	5	17.193	60.855	43.843	1.00	15.57	6
29	OG	SER	A	5	19.340	60.530	41.476	1.00	13.17	8
30	N	VAL	A	6	20.589	59.751	41.567	1.00	10.07	7
31	CA	VAL	A	6	21.169	59.955	42.963	1.00	10.13	6
32	C	VAL	A	6	22.349	59.685	43.172	1.00	10.99	6
33	O	VAL	A	6	21.639	60.160	40.513	1.00	10.81	8
34	CB	VAL	A	6	21.002	59.694	39.148	1.00	13.85	6
35	CG1	VAL	A	6	21.874	61.656	40.459	1.00	15.29	6
36	CG2	VAL	A	6	20.369	60.349	43.964	1.00	12.12	6
37	N	LYS	A	7	20.901	60.604	45.331	1.00	10.30	7
38	CA	LYS	A	7	21.508	59.360	46.015	1.00	9.78	6
39	C	LYS	A	7	22.382	59.560	46.857	1.00	12.06	6
40	O	LYS	A	7	22.963	59.578	46.264	1.00	12.59	8
41	CB	LYS	A	7	19.830	61.187	45.680	1.00	11.40	6
42	CG	LYS	A	7	19.414	62.588	44.966	1.00	12.09	6
43	CD	LYS	A	7	18.160	63.123	46.350	1.00	9.80	6
44	CE	LYS	A	7	17.698	64.488	45.795	1.00	10.87	6
45	NZ	LYS	A	7	17.114	64.187	44.425	1.00	11.38	7
46	N	GLY	A	8	21.036	58.214	45.577	1.00	13.10	7
47	CA	GLY	A	8	21.604	56.982	46.166	1.00	12.31	6
48	C	GLY	A	8	22.718	56.358	45.340	1.00	14.02	6
49	O	GLY	A	8	23.109	55.205	45.579	1.00	13.36	8
50	N	ASP	A	9	23.133	55.048	44.293	1.00	11.90	7
51	CA	ASP	A	9	24.049	56.447	43.319	1.00	11.74	6
52	C	ASP	A	9	25.478	56.996	43.442	1.00	10.18	6
106	CB	ILE	A	15	37.192	59.653	28.343	1.00	10.73	6
107	CG1	ILE	A	15	37.106	60.975	29.131	1.00	11.86	6
108	CG2	ILE	A	15	37.626	60.014	26.904	1.00	12.56	6
109	CD1	ILE	A	15	36.181	62.091	28.574	1.00	15.42	6
110	N	ILE	A	16	34.926	57.450	26.561	1.00	10.37	7
111	CA	ILE	A	16	34.728	56.178	25.868	1.00	11.03	6
112	C	ILE	A	16	35.990	55.729	25.099	1.00	12.23	6
113	O	ILE	A	16	36.342	54.511	25.184	1.00	11.07	8
114	CB	ILE	A	16	33.578	56.292	24.863	1.00	10.56	6
115	CG1	ILE	A	16	32.240	56.387	25.709	1.00	11.92	6
116	CG2	ILE	A	16	33.444	55.053	23.953	1.00	11.18	6
117	CD1	ILE	A	16	31.115	56.958	24.823	1.00	13.67	6
118	N	ASP	A	17	36.565	56.624	24.314	1.00	10.08	7
119	CA	ASP	A	17	37.730	56.165	23.518	1.00	8.61	6
120	C	ASP	A	17	38.911	55.693	24.346	1.00	10.86	6
121	O	ASP	A	17	39.777	54.987	23.831	1.00	11.03	8
122	CB	ASP	A	17	38.184	57.422	22.675	1.00	11.30	6
123	CG	ASP	A	17	39.380	57.017	21.755	1.00	9.77	6
124	OD1	ASP	A	17	39.105	56.206	20.852	1.00	11.65	8
125	OD2	ASP	A	17	40.480	57.562	21.970	1.00	11.48	8
126	N	ARG	A	18	38.972	55.999	25.646	1.00	9.54	7
127	CA	ARG	A	18	40.113	55.719	26.527	1.00	8.38	6
128	C	ARG	A	18	39.826	56.527	27.608	1.00	9.97	6
129	O	ARG	A	18	40.643	54.720	28.501	1.00	13.32	8
130	CB	ARG	A	18	40.537	54.490	27.137	1.00	11.02	6
131	CG	ARG	A	18	40.931	57.083	26.063	1.00	9.63	6
132	CD	ARG	A	18	42.135	58.139	25.237	1.00	9.86	6
133	NE	ARG	A	18	42.280	57.721	23.969	1.00	10.16	7
134	CZ	ARG	A	18	43.748	58.523	23.903	1.00	13.46	6
135	NH1	ARG	A	18	43.350	59.578	24.966	1.00	12.03	7
136	NH2	ARG	A	18	44.350	60.063	22.725	1.00	10.43	7
137	N	PHE	A	19	38.648	54.007	27.497	1.00	11.25	7

138	CA	PHE	A	19	38.296	53.057	28.601	1.00	10.13	6
139	C	PHE	A	19	38.543	51.614	28.192	1.00	12.61	6
140	O	PHE	A	19	39.528	51.024	28.677	1.00	12.51	8
141	CB	PHE	A	19	36.798	53.294	28.945	1.00	13.20	6
142	CG	PHE	A	19	36.342	52.543	30.189	1.00	12.09	6
143	CD1	PHE	A	19	36.849	52.908	31.423	1.00	12.96	6
144	CD2	PHE	A	19	35.472	51.447	30.058	1.00	13.30	6
145	CE1	PHE	A	19	36.500	52.187	32.563	1.00	16.46	6
146	CE2	PHE	A	19	35.184	50.719	31.215	1.00	12.02	6
147	CZ	PHE	A	19	35.638	51.089	32.482	1.00	11.98	6
148	N	TYR	A	20	37.844	51.054	27.199	1.00	11.48	7
149	CA	TYR	A	20	38.154	49.694	26.772	1.00	11.40	6
150	C	TYR	A	20	37.730	49.476	25.321	1.00	10.89	6
151	O	TYR	A	20	36.593	49.813	24.934	1.00	11.80	8
152	CB	TYR	A	20	37.417	48.696	27.719	1.00	12.88	6
153	CG	TYR	A	20	37.927	47.270	27.504	1.00	13.88	6
154	CD1	TYR	A	20	39.216	46.968	27.943	1.00	14.05	6
155	CD2	TYR	A	20	37.160	46.285	26.932	1.00	15.80	6
156	CE1	TYR	A	20	39.717	45.678	27.800	1.00	16.74	7
157	CE2	TYR	A	20	37.658	44.982	26.795	1.00	19.32	6
158	CZ	TYR	A	20	38.935	44.710	27.214	1.00	19.70	6
159	OH	TYR	A	20	39.458	43.422	27.062	1.00	22.69	8
160	N	ASP	A	21	38.662	48.893	24.586	1.00	13.46	7
161	CA	ASP	A	21	38.414	48.628	23.134	1.00	14.35	6
162	C	ASP	A	21	37.754	47.226	23.097	1.00	14.61	6
163	O	ASP	A	21	38.426	46.196	23.063	1.00	13.84	8
164	CB	ASP	A	21	39.746	48.665	22.413	1.00	14.54	6
165	CG	ASP	A	21	39.678	48.444	20.909	1.00	17.30	6
166	OD1	ASP	A	21	38.565	48.288	20.425	1.00	12.67	8
167	OD2	ASP	A	21	40.759	48.450	20.282	1.00	15.87	8
168	N	GLY	A	22	36.430	47.219	23.040	1.00	12.11	7
169	CA	GLY	A	22	35.683	45.934	23.110	1.00	15.21	6
170	C	GLY	A	22	35.482	45.410	21.664	1.00	18.33	6
171	O	GLY	A	22	35.034	44.264	21.516	1.00	17.21	8
172	N	ASP	A	23	35.786	46.189	20.639	1.00	13.30	7
173	CA	ASP	A	23	35.505	45.770	19.261	1.00	14.68	6
174	C	ASP	A	23	36.634	46.389	18.425	1.00	14.54	6
175	O	ASP	A	23	36.570	47.597	18.138	1.00	13.39	8
176	CB	ASP	A	23	34.163	46.271	18.762	1.00	13.59	6
177	CG	ASP	A	23	33.889	45.785	17.319	1.00	16.74	6
178	OD1	ASP	A	23	34.805	45.200	16.750	1.00	17.96	8
179	OD2	ASP	A	23	32.782	46.058	16.872	1.00	16.43	8
180	N	THR	A	24	37.681	45.659	18.083	1.00	15.02	7
181	CA	THR	A	24	38.769	46.308	17.340	1.00	16.37	6
182	C	THR	A	24	38.381	46.606	15.923	1.00	15.39	6
183	O	THR	A	24	39.124	47.376	15.252	1.00	16.40	8
184	CB	THR	A	24	40.033	45.402	17.367	1.00	18.68	6
185	OG1	THR	A	24	39.710	44.147	16.673	1.00	17.99	8
186	CG2	THR	A	24	40.478	45.112	18.772	1.00	22.45	6
187	N	THR	A	25	37.228	46.150	15.423	1.00	16.09	7
188	CA	THR	A	25	36.864	46.330	14.019	1.00	16.07	6
189	C	THR	A	25	36.349	47.743	13.724	1.00	16.52	6
190	O	THR	A	25	36.215	48.013	12.538	1.00	20.25	8
191	CB	THR	A	25	35.780	45.366	13.475	1.00	20.06	6
192	OG1	THR	A	25	34.475	45.532	14.010	1.00	18.04	8
193	CG2	THR	A	25	36.248	43.924	13.739	1.00	21.26	6
194	N	ASN	A	26	36.066	48.509	14.802	1.00	13.98	7
195	CA	ASN	A	26	35.577	49.871	14.471	1.00	12.49	6
196	C	ASN	A	26	36.689	50.862	14.772	1.00	13.05	6
197	O	ASN	A	26	36.435	52.053	14.825	1.00	11.58	8
198	CB	ASN	A	26	34.283	50.103	15.246	1.00	13.85	6
199	CG	ASN	A	26	34.435	49.981	16.739	1.00	15.62	6
200	OD1	ASN	A	26	35.558	50.066	17.224	1.00	12.89	8
201	ND2	ASN	A	26	33.339	49.796	17.497	1.00	16.36	7
202	N	ASN	A	27	37.946	50.401	14.890	1.00	14.79	7
203	CA	ASN	A	27	38.972	51.353	15.290	1.00	12.19	6
204	C	ASN	A	27	39.404	52.337	14.189	1.00	13.62	6
205	O	ASN	A	27	39.775	53.461	14.499	1.00	14.13	8
206	CB	ASN	A	27	40.235	50.575	15.662	1.00	12.10	6
207	CG	ASN	A	27	40.150	49.884	17.001	1.00	15.80	6
208	OD1	ASN	A	27	39.065	49.932	17.554	1.00	12.92	8
209	ND2	ASN	A	27	41.187	49.291	17.571	1.00	15.14	7
210	N	ASN	A	28	39.211	51.954	12.920	1.00	14.26	7
211	CA	ASN	A	28	39.604	52.918	11.854	1.00	15.71	6
212	C	ASN	A	28	38.672	52.705	10.661	1.00	14.39	6
213	O	ASN	A	28	39.059	52.148	9.622	1.00	16.82	8
214	CB	ASN	A	28	41.036	52.497	11.478	1.00	13.79	6
215	CG	ASN	A	28	41.790	53.538	10.656	1.00	21.29	6
216	OD1	ASN	A	28	41.391	54.685	10.535	1.00	17.46	8
217	ND2	ASN	A	28	42.936	53.086	10.108	1.00	24.72	7
218	N	PRO	A	29	37.442	53.154	10.790	1.00	15.04	7
219	CA	PRO	A	29	36.430	52.993	9.742	1.00	17.37	6
220	C	PRO	A	29	36.734	53.802	8.507	1.00	18.08	6
221	O	PRO	A	29	37.259	54.906	8.580	1.00	16.51	8

222	CB	PRO	A	29	35.087	53.483	10.312	1.00	17.71	6
223	CG	PRO	A	29	35.394	53.615	11.787	1.00	17.95	6
224	CD	PRO	A	29	36.907	53.841	11.957	1.00	15.54	6
225	N	ALA	A	30	36.329	53.244	7.331	1.00	16.99	7
226	CA	ALA	A	30	36.533	54.024	6.117	1.00	19.06	6
227	C	ALA	A	30	35.841	55.375	6.161	1.00	16.15	6
228	O	ALA	A	30	36.398	56.355	5.599	1.00	18.27	8
229	CB	ALA	A	30	35.998	53.268	4.880	1.00	21.27	6
230	N	LYS	A	31	34.697	55.514	6.833	1.00	15.01	7
231	CA	LYS	A	31	34.012	56.812	6.886	1.00	14.64	6
232	C	LYS	A	31	34.944	57.908	7.416	1.00	15.07	6
233	O	LYS	A	31	34.722	59.094	7.172	1.00	14.25	8
234	CB	LYS	A	31	32.771	56.667	7.818	1.00	13.99	6
235	CG	LYS	A	31	31.981	57.980	8.050	1.00	13.23	6
236	CD	LYS	A	31	30.617	57.569	8.669	1.00	16.17	6
237	CE	LYS	A	31	29.763	58.766	9.053	1.00	14.65	6
238	NZ	LYS	A	31	30.427	59.568	10.156	1.00	12.05	7
239	N	SER	A	32	35.822	57.610	8.364	1.00	16.27	7
240	CA	SER	A	32	36.675	58.587	9.038	1.00	14.39	6
241	C	SER	A	32	38.087	57.989	9.161	1.00	17.62	6
242	O	SER	A	32	38.770	57.968	10.193	1.00	17.19	8
243	CB	SER	A	32	36.100	58.851	10.460	1.00	12.32	6
244	OG	SER	A	32	35.874	57.664	11.167	1.00	12.92	8
245	N	TYR	A	33	38.596	57.524	8.010	1.00	14.71	7
246	CA	TYR	A	33	39.875	56.801	8.045	1.00	15.23	6
247	C	TYR	A	33	41.051	57.676	8.444	1.00	14.94	6
248	O	TYR	A	33	41.042	58.848	8.023	1.00	17.62	8
249	CB	TYR	A	33	40.075	56.295	6.582	1.00	18.09	6
250	CG	TYR	A	33	41.166	55.254	6.536	1.00	20.46	6
251	CD1	TYR	A	33	40.982	53.978	7.004	1.00	26.95	6
252	CD2	TYR	A	33	42.408	55.618	6.002	1.00	31.14	6
253	CE1	TYR	A	33	41.994	53.027	6.944	1.00	33.56	8
254	CE2	TYR	A	33	43.422	54.670	5.943	1.00	31.30	7
255	CZ	TYR	A	33	43.210	53.409	6.402	1.00	33.88	6
256	OH	TYR	A	33	44.235	52.483	6.334	1.00	44.90	8
257	N	GLY	A	34	42.039	57.105	9.114	1.00	12.92	7
258	CA	GLY	A	34	43.281	57.836	9.403	1.00	14.92	6
259	C	GLY	A	34	43.255	58.672	10.686	1.00	15.08	6
260	O	GLY	A	34	44.274	59.342	10.956	1.00	15.24	8
261	N	LEU	A	35	42.253	58.417	11.548	1.00	12.52	7
262	CA	LEU	A	35	42.215	59.140	12.846	1.00	11.02	6
263	C	LEU	A	35	42.519	58.271	14.028	1.00	15.01	6

264	O	LEU	A	35	42.472	58.755	15.182	1.00	14.18	8
265	CB	LEU	A	35	40.784	59.737	13.090	1.00	10.53	6
266	CG	LEU	A	35	40.170	60.460	11.891	1.00	12.89	6
267	CD1	LEU	A	35	38.783	61.033	12.240	1.00	13.47	6
268	CD2	LEU	A	35	41.090	61.613	11.433	1.00	15.36	6
269	N	TYR	A	36	43.025	57.036	13.757	1.00	13.30	7
270	CA	TYR	A	36	43.335	56.061	14.796	1.00	13.00	6
271	C	TYR	A	36	44.826	55.913	15.032	1.00	13.00	6
272	O	TYR	A	36	45.610	55.855	14.049	1.00	15.18	8
273	CB	TYR	A	36	42.749	54.728	14.291	1.00	15.86	6
274	CG	TYR	A	36	43.149	53.492	15.076	1.00	13.61	6
275	CD1	TYR	A	36	42.927	53.383	16.454	1.00	12.30	6
276	CD2	TYR	A	36	43.807	52.460	14.391	1.00	14.21	6
277	CE1	TYR	A	36	43.317	52.206	17.122	1.00	16.88	6
278	CE2	TYR	A	36	44.182	51.320	15.075	1.00	14.97	6
279	CZ	TYR	A	36	43.930	51.206	16.416	1.00	19.02	6
280	OH	TYR	A	36	44.299	50.063	17.135	1.00	17.90	8
281	N	ASP	A	37	45.211	55.848	16.289	1.00	18.88	7
282	CA	ASP	A	37	46.646	55.624	16.621	1.00	12.45	6
283	C	ASP	A	37	46.700	54.350	17.441	1.00	12.80	6
284	O	ASP	A	37	46.507	54.281	18.673	1.00	14.03	8
285	CB	ASP	A	37	47.120	56.834	17.463	1.00	13.32	6
286	CG	ASP	A	37	48.543	56.543	17.991	1.00	13.46	6
287	OD1	ASP	A	37	49.278	55.720	17.366	1.00	20.57	8
288	OD2	ASP	A	37	48.902	57.113	19.028	1.00	17.00	8
289	N	PRO	A	38	47.163	53.245	16.821	1.00	17.32	7
290	CA	PRO	A	38	47.375	52.024	17.548	1.00	15.62	6
291	C	PRO	A	38	48.484	52.056	18.558	1.00	15.67	6
292	O	PRO	A	38	48.513	51.189	19.436	1.00	15.69	8
293	CB	PRO	A	38	47.669	50.946	16.450	1.00	18.95	6
294	CG	PRO	A	38	48.367	51.843	15.437	1.00	17.01	6
295	CD	PRO	A	38	47.570	53.192	15.409	1.00	18.68	6
296	N	THR	A	39	49.385	53.031	18.514	1.00	18.22	7
297	CA	THR	A	39	50.469	53.080	19.499	1.00	15.60	6
298	C	THR	A	39	50.126	53.773	20.822	1.00	14.85	6
299	O	THR	A	39	50.961	53.777	21.719	1.00	17.68	8
300	CB	THR	A	39	51.692	53.847	18.947	1.00	17.01	6
301	OG1	THR	A	39	51.503	53.239	18.723	1.00	19.14	6
302	CG2	THR	A	39	52.083	53.233	17.573	1.00	16.66	8
303	N	LYS	A	40	48.983	54.487	20.832	1.00	14.93	7
304	CA	LYS	A	40	48.588	55.225	22.041	1.00	14.22	6
305	C	LYS	A	40	49.736	56.141	22.483	1.00	17.80	6

#	Atom	Res	Chain	ResNum	X	Y	Z	Occ	B	El
306	O	LYS	A	40	50.009	56.348	23.685	1.00	17.78	8
307	CB	LYS	A	40	48.104	54.324	23.207	1.00	19.03	6
308	CG	LYS	A	40	47.023	53.320	22.775	1.00	18.65	6
309	CD	LYS	A	40	46.535	52.543	24.031	1.00	21.38	6
310	CE	LYS	A	40	45.432	51.573	23.590	1.00	22.34	6
311	NZ	LYS	A	40	45.883	50.563	22.605	1.00	21.85	7
312	N	SER	A	41	50.307	56.831	21.475	1.00	16.33	7
313	CA	SER	A	41	51.307	57.853	21.746	1.00	17.21	6
314	C	SER	A	41	50.929	59.210	21.203	1.00	16.87	6
315	O	SER	A	41	51.606	60.250	21.492	1.00	16.04	8
316	CB	SER	A	41	52.714	57.429	21.198	1.00	17.96	6
317	OG	SER	A	41	52.625	57.387	19.782	1.00	20.42	8
318	N	LYS	A	42	49.895	59.315	20.388	1.00	13.73	7
319	CA	LYS	A	42	49.446	60.589	19.836	1.00	12.77	6
320	C	LYS	A	42	48.152	60.921	20.603	1.00	13.20	6
321	O	LYS	A	42	47.111	60.351	20.317	1.00	12.98	8
322	CB	LYS	A	42	49.193	60.477	18.321	1.00	14.60	6
323	CG	LYS	A	42	50.523	60.079	17.606	1.00	19.41	6
324	CD	LYS	A	42	50.228	60.163	16.078	1.00	25.03	6
325	CE	LYS	A	42	51.611	60.340	15.395	1.00	34.65	6
326	NZ	LYS	A	42	52.071	58.949	15.130	1.00	41.02	7
327	N	TRP	A	43	48.256	61.858	21.565	1.00	11.08	7
328	CA	TRP	A	43	47.235	61.925	22.643	1.00	13.35	6
329	C	TRP	A	43	45.915	62.494	22.162	1.00	11.08	6
330	O	TRP	A	43	45.002	62.429	22.997	1.00	13.47	8
331	CB	TRP	A	43	47.831	62.848	23.743	1.00	14.15	6
332	CG	TRP	A	43	48.739	61.957	24.592	1.00	12.91	6
333	CD1	TRP	A	43	50.014	61.590	24.338	1.00	14.88	6
334	CD2	TRP	A	43	48.362	61.357	25.845	1.00	12.18	6
335	NE1	TRP	A	43	50.507	60.770	25.364	1.00	16.61	7
336	CE2	TRP	A	43	49.467	60.633	26.297	1.00	17.08	6
337	CE3	TRP	A	43	47.186	61.367	26.617	1.00	13.97	6
338	CZ2	TRP	A	43	49.497	59.891	27.501	1.00	19.44	6
339	CZ3	TRP	A	43	47.223	60.644	27.814	1.00	14.34	6
340	CH2	TRP	A	43	48.333	59.925	28.265	1.00	15.92	6
341	N	LYS	A	44	45.846	63.088	20.972	1.00	11.78	7
342	CA	LYS	A	44	44.532	63.606	20.529	1.00	10.59	6
343	C	LYS	A	44	43.959	62.797	19.362	1.00	11.15	6
344	O	LYS	A	44	43.021	63.227	18.707	1.00	11.48	8
345	CB	LYS	A	44	44.647	65.112	20.097	1.00	11.58	6
346	CG	LYS	A	44	45.053	65.911	21.382	1.00	11.48	6
347	CD	LYS	A	44	44.928	67.435	21.011	1.00	12.19	6
348	CE	LYS	A	44	45.254	68.171	22.334	1.00	15.83	6
349	NZ	LYS	A	44	45.125	69.681	22.068	1.00	18.92	7
350	N	MET	A	45	44.473	61.576	19.114	1.00	10.42	7
351	CA	MET	A	45	43.881	60.686	18.112	1.00	12.24	6
352	C	MET	A	45	42.952	59.664	18.768	1.00	11.36	6
353	O	MET	A	45	43.011	59.512	19.985	1.00	12.88	8
354	CB	MET	A	45	45.028	59.874	17.442	1.00	13.26	6
355	CG	AMET	A	45	46.067	60.710	16.692	0.50	14.78	6
356	SD	AMET	A	45	45.379	61.237	15.135	0.50	13.95	16
357	CE	AMET	A	45	45.728	60.040	13.903	0.50	12.41	6
355	CG	BMET	A	45	45.776	60.960	16.619	0.50	11.59	6
356	SD	BMET	A	45	46.918	60.290	15.431	0.50	16.20	16
357	CE	BMET	A	45	45.864	59.453	14.271	0.50	18.11	6
358	N	TYR	A	46	42.122	58.961	17.976	1.00	10.91	7
359	CA	TYR	A	46	41.356	57.880	18.584	1.00	13.29	6
360	C	TYR	A	46	42.263	56.691	18.938	1.00	13.10	6
361	O	TYR	A	46	43.076	56.318	18.094	1.00	12.46	8
362	CB	TYR	A	46	40.258	57.364	17.660	1.00	12.44	6
363	CG	TYR	A	46	39.031	58.210	17.416	1.00	13.02	6
364	CD1	TYR	A	46	39.075	59.210	16.436	1.00	11.30	6
365	CD2	TYR	A	46	37.846	57.978	18.105	1.00	12.45	6
366	CE1	TYR	A	46	37.940	59.997	16.146	1.00	12.75	6
367	CE2	TYR	A	46	36.683	58.746	17.838	1.00	9.77	6
368	CZ	TYR	A	46	36.789	59.707	16.881	1.00	10.60	6
369	OH	TYR	A	46	35.703	60.490	16.547	1.00	11.65	8
370	N	TRP	A	47	42.097	56.222	20.188	1.00	9.67	7
371	CA	TRP	A	47	42.866	55.089	20.664	1.00	11.50	6
372	C	TRP	A	47	42.065	53.770	20.579	1.00	12.29	6
373	O	TRP	A	47	42.633	52.676	20.711	1.00	12.20	8
374	CB	TRP	A	47	43.430	55.285	22.077	1.00	12.80	6
375	CG	TRP	A	47	44.548	56.316	22.086	1.00	10.46	6
376	CD1	TRP	A	47	45.068	57.007	21.037	1.00	11.88	6
377	CD2	TRP	A	47	45.300	56.687	23.218	1.00	10.01	6
378	NE1	TRP	A	47	46.060	57.853	21.485	1.00	11.36	7
379	CE2	TRP	A	47	46.219	57.700	22.820	1.00	12.07	6
380	CE3	TRP	A	47	45.198	56.392	24.603	1.00	12.09	7
381	CZ2	TRP	A	47	47.103	58.301	23.715	1.00	13.05	6
382	CZ3	TRP	A	47	46.072	56.974	25.484	1.00	15.07	6
383	CH2	TRP	A	47	47.002	57.939	25.033	1.00	16.33	6
384	N	GLY	A	48	40.752	53.875	20.442	1.00	10.96	7
385	CA	GLY	A	48	39.995	52.631	20.097	1.00	11.53	6
386	C	GLY	A	48	38.960	52.197	21.106	1.00	11.03	6

387	O	GLY	A	48	38.208	51.215	20.845	1.00	12.01	8
388	N	GLY	A	49	38.834	52.862	22.221	1.00	12.42	7
389	CA	GLY	A	49	37.789	52.443	23.230	1.00	12.08	6
390	C	GLY	A	49	36.451	52.679	22.614	1.00	9.81	6
391	O	GLY	A	49	36.173	53.629	21.880	1.00	10.92	8
392	N	ASP	A	50	35.433	51.851	23.065	1.00	10.42	7
393	CA	ASP	A	50	34.135	51.985	22.429	1.00	11.91	6
394	C	ASP	A	50	32.977	51.516	23.344	1.00	11.85	6
395	O	ASP	A	50	33.188	51.228	24.489	1.00	12.76	8
396	CB	ASP	A	50	34.148	51.188	21.094	1.00	10.66	6
397	CG	ASP	A	50	34.693	49.790	21.327	1.00	14.50	6
398	OD1	ASP	A	50	34.446	49.184	22.384	1.00	11.19	8
399	OD2	ASP	A	50	35.425	49.205	20.532	1.00	11.87	8
400	N	LEU	A	51	31.762	51.615	22.778	1.00	11.80	7
401	CA	LEU	A	51	30.580	51.320	23.617	1.00	11.32	6
402	C	LEU	A	51	30.568	49.843	23.973	1.00	13.43	6
403	O	LEU	A	51	30.145	49.499	25.090	1.00	11.80	8
404	CB	LEU	A	51	29.272	51.662	22.869	1.00	12.03	6
405	CG	LEU	A	51	29.178	53.205	22.638	1.00	11.71	6
406	CD1	LEU	A	51	28.036	53.389	21.666	1.00	13.88	6
407	CD2	LEU	A	51	28.915	53.930	23.954	1.00	15.76	7
408	N	GLU	A	52	30.942	48.987	23.037	1.00	12.67	6
409	CA	GLU	A	52	30.995	47.541	23.443	1.00	12.25	6
410	C	GLU	A	52	32.024	47.239	24.516	1.00	12.73	6
411	O	GLU	A	52	31.816	46.375	25.382	1.00	13.32	8
412	CB	GLU	A	52	31.182	46.786	22.122	1.00	16.82	6
413	CG	GLU	A	52	31.390	45.298	22.295	1.00	22.57	6
414	CD	GLU	A	52	30.227	44.545	22.992	1.00	12.69	6
415	OE1	GLU	A	52	29.097	45.029	23.005	1.00	17.98	8
416	OE2	GLU	A	52	30.680	43.475	23.419	1.00	16.49	8
417	N	GLY	A	53	33.114	48.012	24.628	1.00	12.03	7
418	CA	GLY	A	53	34.108	47.857	25.680	1.00	13.18	6
419	C	GLY	A	53	33.471	48.292	27.005	1.00	12.67	6
420	O	GLY	A	53	33.737	47.586	28.000	1.00	11.91	8
421	N	VAL	A	54	32.653	49.355	27.005	1.00	11.80	7
422	CA	VAL	A	54	31.996	49.680	28.280	1.00	10.05	6
423	C	VAL	A	54	31.078	48.502	28.715	1.00	12.37	6
424	O	VAL	A	54	31.055	48.111	29.879	1.00	12.15	8
425	CB	VAL	A	54	31.154	50.947	28.220	1.00	11.03	6
426	CG1	VAL	A	54	30.449	51.255	29.552	1.00	13.86	6
427	CG2	VAL	A	54	32.100	52.143	27.853	1.00	11.86	6
428	N	ARG	A	55	30.387	47.952	27.708	1.00	9.95	7
429	CA	ARG	A	55	29.382	46.875	28.101	1.00	13.29	6
430	C	ARG	A	55	30.112	45.671	28.652	1.00	12.91	6
431	O	ARG	A	55	29.684	44.943	29.596	1.00	13.96	8
432	CB	ARG	A	55	28.627	46.458	26.819	1.00	13.43	6
433	CG	ARG	A	55	27.364	45.611	27.165	1.00	13.64	6
434	CD	ARG	A	55	26.723	44.974	25.877	1.00	13.15	6
435	NE	ARG	A	55	27.745	44.040	25.358	1.00	13.30	7
436	CZ	ARG	A	55	28.117	42.905	25.921	1.00	14.35	6
437	NH1	ARG	A	55	27.475	42.404	27.011	1.00	15.82	7
438	NH2	ARG	A	55	29.125	42.171	25.446	1.00	17.12	7
439	N	GLN	A	56	31.265	45.354	28.031	1.00	11.75	7
440	CA	GLN	A	56	32.050	44.171	28.503	1.00	12.69	6
441	C	GLN	A	56	32.530	44.339	29.945	1.00	14.76	6
442	O	GLN	A	56	32.895	43.338	30.611	1.00	15.16	8
443	CB	GLN	A	56	33.249	43.948	27.536	1.00	12.12	6
444	CG	GLN	A	56	32.718	43.310	26.223	1.00	12.45	6
445	CD	GLN	A	56	33.748	43.189	25.110	1.00	18.74	6
446	OE1	GLN	A	56	33.441	43.161	23.879	1.00	21.22	8
447	NE2	GLN	A	56	34.957	43.066	25.540	1.00	13.29	7
448	N	LYS	A	57	32.816	45.574	30.355	1.00	13.93	7
449	CA	LYS	A	57	33.243	45.881	31.703	1.00	11.77	6
450	C	LYS	A	57	32.146	46.200	32.702	1.00	13.05	6
451	O	LYS	A	57	32.397	46.651	33.834	1.00	12.11	8
452	CB	LYS	A	57	34.240	47.112	31.625	1.00	12.27	6
453	CG	LYS	A	57	35.508	46.752	30.818	1.00	13.15	6
454	CD	LYS	A	57	36.167	45.442	31.318	1.00	13.38	6
455	CE	LYS	A	57	37.577	45.277	30.729	1.00	16.88	6
456	NZ	LYS	A	57	38.170	43.960	31.261	1.00	17.21	7
457	N	LEU	A	58	30.883	45.891	32.388	1.00	12.90	7
458	CA	LEU	A	58	29.789	46.048	33.338	1.00	14.31	6
459	C	LEU	A	58	29.981	45.299	34.668	1.00	12.68	6
460	O	LEU	A	58	29.737	45.732	35.732	1.00	13.94	8
461	CB	LEU	A	58	28.407	45.779	32.723	1.00	12.52	6
462	CG	LEU	A	58	27.963	46.878	31.718	1.00	12.14	6
463	CD1	LEU	A	58	26.709	46.366	30.943	1.00	14.87	6
464	CD2	LEU	A	58	27.586	48.136	32.488	1.00	15.84	6
465	N	PRO	A	59	30.555	44.107	34.670	1.00	13.13	7
466	CA	PRO	A	59	30.776	43.396	35.937	1.00	14.64	6
467	C	PRO	A	59	31.759	44.139	36.827	1.00	14.63	6
468	O	PRO	A	59	31.532	44.250	38.038	1.00	15.79	8
469	CB	PRO	A	59	31.436	42.034	35.525	1.00	15.40	6
470	CG	PRO	A	59	30.719	41.845	34.161	1.00	16.79	6

#	Atom	Res	Chain	Res#	x	y	z	Occ	B	El
471	CD	PRO	A	59	30.807	43.247	33.514	1.00	16.71	6
472	N	TYR	A	60	32.806	44.717	36.210	1.00	12.89	7
473	CA	TYR	A	60	33.789	45.511	36.994	1.00	12.47	6
474	C	TYR	A	60	33.072	46.731	37.584	1.00	12.64	6
475	O	TYR	A	60	33.237	46.994	38.797	1.00	13.63	8
476	CB	TYR	A	60	34.918	45.920	36.026	1.00	12.16	6
477	CG	TYR	A	60	35.856	46.938	36.667	1.00	12.17	6
478	CD1	TYR	A	60	36.917	46.528	37.462	1.00	13.23	6
479	CD2	TYR	A	60	35.602	48.293	36.453	1.00	12.24	6
480	CE1	TYR	A	60	37.730	47.509	38.049	1.00	12.99	6
481	CE2	TYR	A	60	36.438	49.268	37.022	1.00	14.93	6
482	CZ	TYR	A	60	37.473	48.852	37.823	1.00	14.75	6
483	OH	TYR	A	60	38.287	49.782	38.464	1.00	13.93	6
484	N	LEU	A	61	32.298	47.410	36.735	1.00	11.74	6
485	CA	LEU	A	61	31.622	48.610	37.225	1.00	11.91	6
486	C	LEU	A	61	30.570	48.316	38.272	1.00	14.11	6
487	O	LEU	A	61	30.508	49.022	39.283	1.00	13.33	8
488	CB	LEU	A	61	30.993	49.382	36.051	1.00	12.06	6
489	CG	LEU	A	61	32.030	49.809	34.992	1.00	13.18	6
490	CD1	LEU	A	61	31.263	50.310	33.753	1.00	15.35	6
491	CD2	LEU	A	61	32.865	50.971	35.605	1.00	16.34	6
492	N	LYS	A	62	29.850	47.217	38.162	1.00	12.47	7
493	CA	LYS	A	62	28.890	46.844	39.202	1.00	13.25	6
494	C	LYS	A	62	29.614	46.558	40.535	1.00	13.60	6
495	O	LYS	A	62	29.149	47.032	41.576	1.00	15.97	8
496	CB	LYS	A	62	28.117	45.588	38.730	1.00	14.71	6
497	CG	LYS	A	62	27.011	45.263	39.764	1.00	17.27	6
498	CD	LYS	A	62	25.908	44.363	39.223	1.00	30.17	6
499	CE	LYS	A	62	24.879	44.088	40.343	1.00	25.52	6
500	NZ	LYS	A	62	23.887	45.203	40.515	1.00	24.63	7
501	N	GLN	A	63	30.722	45.781	40.455	1.00	12.05	7
502	CA	GLN	A	63	31.437	45.448	41.660	1.00	10.80	6
503	C	GLN	A	63	32.010	46.705	42.325	1.00	13.71	6
504	O	GLN	A	63	32.200	46.751	43.544	1.00	13.66	8
505	CB	AGLN	A	63	32.582	44.501	41.262	0.66	17.93	6
506	CG	AGLN	A	63	32.178	43.092	40.865	0.66	27.01	6
507	CD	AGLN	A	63	33.421	42.362	40.343	0.66	35.95	6
508	OE1	AGLN	A	63	34.283	42.022	41.149	0.66	40.99	8
509	NE2	AGLN	A	63	33.525	42.128	39.043	0.66	36.04	7
505	CB	BGLN	A	63	32.511	44.387	41.345	0.33	7.59	6
506	CG	BGLN	A	63	33.072	43.793	42.621	0.33	8.29	6
507	CD	BGLN	A	63	34.234	42.824	42.408	0.33	8.28	6

#	Atom	Res	Chain	Res#	x	y	z	Occ	B	El
508	OE1	BGLN	A	63	34.480	42.416	41.279	0.33	14.24	8
509	NE2	BGLN	A	63	34.884	42.519	43.537	0.33	12.44	7
510	N	LEU	A	64	32.481	47.646	41.498	1.00	10.99	7
511	CA	LEU	A	64	32.993	48.909	42.087	1.00	15.73	6
512	C	LEU	A	64	31.893	49.656	42.837	1.00	14.21	8
513	O	LEU	A	64	32.253	50.516	43.659	1.00	14.81	6
514	CB	LEU	A	64	33.536	49.777	40.930	1.00	14.15	6
515	CG	LEU	A	64	34.050	51.201	41.274	1.00	13.04	6
516	CD1	LEU	A	64	35.177	51.132	42.303	1.00	12.46	6
517	CD2	LEU	A	64	34.587	51.825	39.963	1.00	12.70	6
518	N	GLY	A	65	30.605	49.492	42.566	1.00	14.23	7
519	CA	GLY	A	65	29.537	50.247	43.205	1.00	13.78	6
520	C	GLY	A	65	28.987	51.337	42.311	1.00	14.56	6
521	O	GLY	A	65	28.207	52.222	42.758	1.00	13.13	8
522	N	VAL	A	66	29.343	51.265	41.014	1.00	12.13	7
523	CA	VAL	A	66	28.773	52.267	40.114	1.00	10.77	6
524	C	VAL	A	66	27.297	52.007	39.842	1.00	13.82	6
525	O	VAL	A	66	26.933	50.836	39.617	1.00	13.38	8
526	CB	VAL	A	66	29.491	52.192	38.744	1.00	11.10	6
527	CG1	VAL	A	66	28.892	53.220	37.731	1.00	12.12	6
528	CG2	VAL	A	66	30.961	52.489	38.974	1.00	14.48	6
529	N	THR	A	67	26.431	53.016	39.992	1.00	10.70	7
530	CA	THR	A	67	25.022	52.822	39.675	1.00	12.22	6
531	C	THR	A	67	24.526	53.737	38.565	1.00	13.77	6
532	O	THR	A	67	23.404	53.538	38.103	1.00	13.29	8
533	CB	THR	A	67	24.072	52.926	40.898	1.00	14.25	6
534	OG1	THR	A	67	24.680	53.791	41.874	1.00	13.89	6
535	CG2	THR	A	67	24.085	51.519	41.584	1.00	14.96	8
536	N	THR	A	68	25.351	54.723	38.156	1.00	12.63	6
537	CA	THR	A	68	25.042	55.479	36.914	1.00	10.68	7
538	C	THR	A	68	26.379	55.684	36.193	1.00	9.37	7
539	O	THR	A	68	27.313	56.214	36.775	1.00	11.86	6
540	CB	THR	A	68	24.388	56.837	37.236	1.00	13.89	6
541	OG1	THR	A	68	23.094	56.649	37.793	1.00	13.19	8
542	CG2	THR	A	68	24.269	57.734	35.965	1.00	14.32	6
543	N	ILE	A	69	26.394	55.287	34.896	1.00	9.42	6
544	CA	ILE	A	69	27.605	55.600	34.073	1.00	8.17	7
545	C	ILE	A	69	27.249	56.897	33.377	1.00	11.16	6
546	O	ILE	A	69	26.200	57.080	32.734	1.00	12.50	8
547	CB	ILE	A	69	27.714	54.510	32.986	1.00	12.66	6
548	CG1	ILE	A	69	28.160	53.214	33.736	1.00	14.05	6
549	CG2	ILE	A	69	28.737	54.893	31.900	1.00	11.22	6

Serial	Atom	Res	Chain	ResNum	X	Y	Z	Occ	B	El
550	CD1	ILE	A	69	27.937	52.008	32.775	1.00	13.93	6
551	N	TRP	A	70	28.196	57.872	33.425	1.00	9.74	7
552	CA	TRP	A	70	28.095	59.093	32.600	1.00	10.34	6
553	C	TRP	A	70	28.991	58.790	31.403	1.00	11.63	6
554	O	TRP	A	70	30.214	58.751	31.558	1.00	11.51	8
555	CB	TRP	A	70	28.494	60.327	33.441	1.00	9.83	6
556	CG	TRP	A	70	28.954	61.558	32.738	1.00	8.81	6
557	CD1	TRP	A	70	29.050	61.770	31.360	1.00	13.03	6
558	CD2	TRP	A	70	29.591	62.698	33.356	1.00	10.57	6
559	NE1	TRP	A	70	29.645	63.016	31.118	1.00	12.41	7
560	CE2	TRP	A	70	30.017	63.558	32.338	1.00	10.48	6
561	CE3	TRP	A	70	29.830	63.007	34.699	1.00	12.18	6
562	CZ2	TRP	A	70	30.721	64.729	32.587	1.00	9.61	6
563	CZ3	TRP	A	70	30.426	64.234	34.950	1.00	11.62	6
564	CH2	TRP	A	70	30.896	65.061	33.914	1.00	13.98	6
565	N	LEU	A	71	28.373	58.542	30.225	1.00	10.83	7
566	CA	LEU	A	71	29.219	58.311	29.023	1.00	11.92	6
567	C	LEU	A	71	29.585	59.691	28.439	1.00	10.49	6
568	O	LEU	A	71	28.669	60.552	28.276	1.00	10.64	8
569	CB	LEU	A	71	28.342	57.617	27.923	1.00	11.13	6
570	CG	LEU	A	71	27.991	56.159	28.240	1.00	11.31	6
571	CD1	LEU	A	71	27.073	55.665	27.096	1.00	10.98	6
572	CD2	LEU	A	71	29.253	55.314	28.322	1.00	11.96	6
573	N	SER	A	72	30.870	59.865	28.066	1.00	10.17	7
574	CA	SER	A	72	31.250	60.995	27.218	1.00	9.81	6
575	C	SER	A	72	30.455	60.988	25.920	1.00	11.22	6
576	O	SER	A	72	29.733	60.011	25.572	1.00	10.33	8
577	CB	SER	A	72	32.773	60.898	26.944	1.00	10.62	6
578	OG	SER	A	72	33.092	59.694	26.237	1.00	11.62	8
579	N	PRO	A	73	30.447	62.034	25.128	1.00	11.10	7
580	CA	PRO	A	73	29.427	62.188	24.048	1.00	11.86	6
581	C	PRO	A	73	29.521	61.057	23.042	1.00	13.54	6
582	O	PRO	A	73	30.653	60.649	22.674	1.00	11.75	8
583	CB	PRO	A	73	29.672	63.563	23.414	1.00	10.32	6
584	CG	PRO	A	73	30.360	64.313	24.557	1.00	10.23	6
585	CD	PRO	A	73	31.228	63.286	25.358	1.00	11.09	6
586	N	VAL	A	74	28.345	60.538	22.623	1.00	11.02	7
587	CA	VAL	A	74	28.351	59.338	21.794	1.00	9.14	6
588	C	VAL	A	74	27.998	59.628	20.344	1.00	9.79	6
589	O	VAL	A	74	28.041	58.700	19.549	1.00	10.96	8
590	CB	VAL	A	74	27.260	58.313	22.311	1.00	9.29	6
591	CG1	VAL	A	74	27.541	57.935	23.780	1.00	11.66	6
592	CG2	VAL	A	74	25.882	58.894	22.161	1.00	10.26	6
593	N	LEU	A	75	27.708	60.912	19.980	1.00	10.62	7
594	CA	LEU	A	75	27.182	61.136	18.619	1.00	10.48	6
595	C	LEU	A	75	28.305	61.464	17.620	1.00	12.02	6
596	O	LEU	A	75	29.436	61.678	18.016	1.00	11.28	8
597	CB	LEU	A	75	26.111	62.279	18.660	1.00	10.81	6
598	CG	LEU	A	75	24.952	61.966	19.634	1.00	11.49	6
599	CD1	LEU	A	75	24.074	63.178	19.911	1.00	11.47	6
600	CD2	LEU	A	75	24.074	60.864	18.960	1.00	11.05	6
601	N	ASP	A	76	27.958	61.296	16.347	1.00	10.11	7
602	CA	ASP	A	76	29.020	61.292	15.299	1.00	9.23	6
603	C	ASP	A	76	29.821	62.605	15.351	1.00	9.60	6
604	O	ASP	A	76	29.263	63.683	15.155	1.00	11.11	8
605	CB	ASP	A	76	28.264	61.153	13.979	1.00	9.89	6
606	CG	ASP	A	76	29.177	61.079	12.745	1.00	12.56	6
607	OD1	ASP	A	76	30.380	60.856	12.895	1.00	12.97	8
608	OD2	ASP	A	76	28.617	61.239	11.641	1.00	11.97	8
609	N	ASN	A	77	31.131	62.435	15.524	1.00	9.80	7
610	CA	ASN	A	77	32.043	63.570	15.534	1.00	9.86	6
611	C	ASN	A	77	32.766	63.691	14.180	1.00	10.69	6
612	O	ASN	A	77	32.797	62.759	13.385	1.00	10.51	8
613	CB	ASN	A	77	33.117	63.370	16.619	1.00	10.13	6
614	CG	ASN	A	77	32.685	63.988	17.945	1.00	13.40	6
615	OD1	ASN	A	77	33.515	64.636	18.600	1.00	10.92	8
616	ND2	ASN	A	77	31.412	63.908	18.341	1.00	11.73	7
617	N	LEU	A	78	33.296	64.886	13.967	1.00	10.62	7
618	CA	LEU	A	78	34.240	65.195	12.866	1.00	11.24	6
619	C	LEU	A	78	34.929	63.977	12.309	1.00	8.87	6
620	O	LEU	A	78	35.632	63.257	13.026	1.00	12.04	8
621	CB	LEU	A	78	35.257	66.197	13.506	1.00	9.81	6
622	CG	LEU	A	78	36.289	66.679	12.399	1.00	9.30	6
623	CD1	LEU	A	78	35.622	67.597	11.418	1.00	11.34	6
624	CD2	LEU	A	78	37.382	67.439	13.176	1.00	13.04	6
625	N	ASP	A	79	34.801	63.867	10.945	1.00	11.36	7
626	CA	ASP	A	79	35.393	62.670	10.348	1.00	9.32	6
627	C	ASP	A	79	36.754	62.947	9.688	1.00	12.70	6
628	O	ASP	A	79	37.275	62.042	9.026	1.00	14.30	8
629	CB	ASP	A	79	34.468	62.189	9.168	1.00	14.19	6
630	CG	ASP	A	79	33.217	61.518	9.658	1.00	15.14	6
631	OD1	ASP	A	79	33.208	61.150	10.841	1.00	12.50	8
632	OD2	ASP	A	79	32.239	61.307	8.931	1.00	12.26	8
633	N	THR	A	80	37.307	64.115	9.950	1.00	12.57	7

#	Atom	Res	Chain	ResNum	X	Y	Z	Occ	B	Elem
634	CA	THR	A	80	38.652	64.479	9.456	1.00	14.13	6
635	C	THR	A	80	39.521	64.930	10.635	1.00	13.16	6
636	O	THR	A	80	39.072	64.981	11.769	1.00	12.58	8
637	CB	THR	A	80	38.583	65.712	8.534	1.00	14.19	6
638	OG1	THR	A	80	38.265	66.927	9.264	1.00	15.46	8
639	CG2	THR	A	80	37.593	65.645	7.372	1.00	19.78	6
640	N	LEU	A	81	40.809	65.216	10.327	1.00	13.37	7
641	CA	LEU	A	81	41.651	65.864	11.341	1.00	10.22	6
642	C	LEU	A	81	41.263	67.297	11.488	1.00	12.10	6
643	O	LEU	A	81	40.635	67.900	10.595	1.00	12.72	8
644	CB	LEU	A	81	43.143	65.818	10.830	1.00	10.42	6
645	CG	LEU	A	81	43.643	64.345	10.809	1.00	15.27	6
646	CD1	LEU	A	81	44.897	64.351	9.927	1.00	21.14	6
647	CD2	LEU	A	81	44.059	63.885	12.244	1.00	14.05	6
648	N	ALA	A	82	41.647	67.888	12.629	1.00	11.74	7
649	CA	ALA	A	82	41.548	69.320	12.798	1.00	13.34	6
650	C	ALA	A	82	42.941	69.801	13.202	1.00	12.91	6
651	O	ALA	A	82	43.208	70.247	14.316	1.00	11.11	8
652	CB	ALA	A	82	40.566	69.586	13.989	1.00	14.40	6
653	N	GLY	A	83	43.811	69.835	12.180	1.00	12.89	7
654	CA	GLY	A	83	45.245	70.145	12.484	1.00	11.52	6
655	C	GLY	A	83	45.960	68.860	12.923	1.00	12.02	6
656	O	GLY	A	83	45.405	67.737	13.060	1.00	11.47	8
657	N	THR	A	84	47.262	68.987	13.230	1.00	12.31	7
658	CA	THR	A	84	48.160	67.879	13.444	1.00	11.91	6
659	C	THR	A	84	47.716	66.883	14.496	1.00	10.74	6
660	O	THR	A	84	47.554	67.213	15.687	1.00	11.33	8
661	CB	THR	A	84	49.570	68.477	13.888	1.00	10.32	6
662	OG1	THR	A	84	49.942	69.432	12.873	1.00	12.17	8
663	CG2	THR	A	84	50.533	67.298	14.074	1.00	13.58	6
664	N	ASP	A	85	47.462	65.652	14.019	1.00	12.06	7
665	CA	ASP	A	85	47.117	64.552	14.933	1.00	10.80	6
666	C	ASP	A	85	45.981	64.902	15.894	1.00	12.31	6
667	O	ASP	A	85	45.986	64.416	17.030	1.00	13.43	8
668	CB	ASP	A	85	48.356	64.077	15.747	1.00	14.44	6
669	CG	ASP	A	85	49.500	63.600	14.831	1.00	22.31	6
670	OD1	ASP	A	85	49.284	63.065	13.744	1.00	14.91	8
671	OD2	ASP	A	85	50.645	63.782	15.275	1.00	23.25	8
672	N	ASN	A	86	45.024	65.712	15.418	1.00	10.53	7
673	CA	ASN	A	86	44.024	66.196	16.401	1.00	10.12	6
674	C	ASN	A	86	42.644	65.784	15.880	1.00	9.75	6
675	O	ASN	A	86	42.241	66.073	14.747	1.00	12.30	8
676	CB	ASN	A	86	44.190	67.732	16.468	1.00	11.52	6
677	CG	ASN	A	86	43.470	68.295	17.677	1.00	13.68	6
678	OD1	ASN	A	86	42.743	67.535	18.341	1.00	14.08	8
679	ND2	ASN	A	86	43.644	69.546	18.063	1.00	15.21	7
680	N	THR	A	87	41.877	65.099	16.705	1.00	10.11	7
681	CA	THR	A	87	40.619	64.459	16.251	1.00	10.70	6
682	C	THR	A	87	39.525	64.643	17.317	1.00	11.84	6
683	O	THR	A	87	39.764	65.021	18.471	1.00	11.11	8
684	CB	THR	A	87	40.745	62.924	16.124	1.00	12.30	6
685	OG1	THR	A	87	40.840	62.298	17.444	1.00	11.34	8
686	CG2	THR	A	87	41.961	62.467	15.375	1.00	11.19	6
687	N	GLY	A	88	38.307	64.357	16.903	1.00	11.28	7
688	CA	GLY	A	88	37.184	64.329	17.887	1.00	10.27	6
689	C	GLY	A	88	37.107	63.063	18.721	1.00	10.54	6
690	O	GLY	A	88	35.954	62.712	19.121	1.00	10.21	8
691	N	TYR	A	89	38.196	62.404	19.087	1.00	10.33	7
692	CA	TYR	A	89	38.134	61.241	19.955	1.00	10.06	6
693	C	TYR	A	89	37.314	61.476	21.204	1.00	11.70	6
694	O	TYR	A	89	36.760	60.489	21.732	1.00	11.42	8
695	CB	TYR	A	89	39.564	60.769	20.316	1.00	9.10	6
696	CG	TYR	A	89	40.152	61.653	21.412	1.00	11.07	6
697	CD1	TYR	A	89	40.732	61.106	22.750	1.00	10.61	6
698	CD2	TYR	A	89	40.058	62.857	21.106	1.00	9.12	6
699	CE1	TYR	A	89	41.243	63.688	22.115	1.00	9.21	6
700	CE2	TYR	A	89	40.505	62.054	23.777	1.00	11.90	6
701	CZ	TYR	A	89	41.098	63.267	23.443	1.00	10.92	6
702	OH	TYR	A	89	41.593	64.126	24.411	1.00	10.82	8
703	N	HIS	A	90	37.283	62.703	21.748	1.00	10.85	7
704	CA	HIS	A	90	36.623	62.983	23.011	1.00	9.54	6
705	C	HIS	A	90	35.095	63.104	22.837	1.00	7.69	6
706	O	HIS	A	90	34.392	63.040	23.856	1.00	9.02	8
707	CB	HIS	A	90	37.178	64.338	23.555	1.00	10.73	6
708	CG	HIS	A	90	37.294	65.403	22.507	1.00	11.48	6
709	ND1	HIS	A	90	36.210	66.035	21.926	1.00	9.73	7
710	CD2	HIS	A	90	38.405	65.898	21.906	1.00	8.93	6
711	CE1	HIS	A	90	36.686	66.903	21.010	1.00	10.94	6
712	NE2	HIS	A	90	37.988	66.862	20.995	1.00	10.18	7
713	N	GLY	A	91	34.616	63.356	21.629	1.00	9.66	7
714	CA	GLY	A	91	33.143	63.393	21.404	1.00	9.79	6
715	C	GLY	A	91	32.505	64.768	21.395	1.00	10.82	6
716	O	GLY	A	91	31.287	64.916	21.102	1.00	11.09	8
717	N	TYR	A	92	33.288	65.840	21.633	1.00	10.15	7

#	Atom	Res	Chain	Seq	X	Y	Z	Occ	B	El
718	CA	TYR	A	92	32.653	67.152	21.753	1.00	10.44	6
719	C	TYR	A	92	32.556	67.935	20.459	1.00	8.85	6
720	O	TYR	A	92	32.132	69.141	20.520	1.00	9.64	8
721	CB	TYR	A	92	33.461	67.977	22.837	1.00	10.08	6
722	CG	TYR	A	92	33.217	67.385	24.219	1.00	10.98	6
723	CD1	TYR	A	92	32.091	67.719	24.974	1.00	10.25	6
724	CD2	TYR	A	92	34.112	66.472	24.730	1.00	10.46	6
725	CE1	TYR	A	92	31.905	67.143	26.247	1.00	10.96	6
726	CE2	TYR	A	92	33.932	65.888	25.983	1.00	14.14	6
727	CZ	TYR	A	92	32.829	66.237	26.715	1.00	11.56	6
728	OH	TYR	A	92	32.648	65.665	27.970	1.00	12.31	8
729	N	TRP	A	93	32.968	67.343	19.345	1.00	10.19	7
730	CA	TRP	A	93	32.914	68.040	18.039	1.00	11.73	6
731	C	TRP	A	93	31.949	67.307	17.103	1.00	9.84	6
732	O	TRP	A	93	32.336	66.443	16.294	1.00	10.83	8
733	CB	TRP	A	93	34.322	68.024	17.415	1.00	12.42	6
734	CG	TRP	A	93	35.359	68.775	18.253	1.00	11.75	6
735	CD1	TRP	A	93	35.181	69.659	19.258	1.00	12.58	6
736	CD2	TRP	A	93	36.786	68.657	18.020	1.00	11.96	6
737	NE1	TRP	A	93	36.448	70.120	19.694	1.00	13.60	7
738	CE2	TRP	A	93	37.397	69.501	18.932	1.00	14.78	6
739	CE3	TRP	A	93	37.559	67.884	17.130	1.00	13.74	6
740	CZ2	TRP	A	93	38.808	69.626	19.040	1.00	15.96	6
741	CZ3	TRP	A	93	38.959	68.021	17.209	1.00	10.95	6
742	CH2	TRP	A	93	39.526	68.880	18.171	1.00	9.91	6
743	N	THR	A	94	30.669	67.534	17.237	1.00	10.45	7
744	CA	THR	A	94	29.603	66.690	16.661	1.00	9.91	6
745	C	THR	A	94	29.244	67.242	15.270	1.00	11.32	6
746	O	THR	A	94	28.854	68.415	15.074	1.00	11.48	8
747	CB	THR	A	94	28.302	66.837	17.495	1.00	9.85	6
748	OG1	THR	A	94	28.643	66.563	18.891	1.00	11.49	8
749	CG2	THR	A	94	27.263	65.739	17.079	1.00	11.23	6
750	N	ARG	A	95	29.315	66.299	14.292	1.00	10.33	7
751	CA	ARG	A	95	28.819	66.623	12.945	1.00	11.51	6
752	C	ARG	A	95	27.382	66.123	12.703	1.00	13.19	6
753	O	ARG	A	95	26.834	65.999	11.700	1.00	12.00	8
754	CB	ARG	A	95	29.766	65.999	11.920	1.00	12.95	6
755	CG	ARG	A	95	29.739	64.437	11.900	1.00	9.03	6
756	CD	ARG	A	95	30.894	63.964	11.008	1.00	12.99	6
757	NE	ARG	A	95	30.917	64.346	9.584	1.00	13.10	7
758	CZ	ARG	A	95	30.216	63.680	8.633	1.00	13.74	6
759	NH1	ARG	A	95	29.307	62.746	8.918	1.00	13.30	7
760	NH2	ARG	A	95	30.477	64.026	7.380	1.00	15.12	7
761	N	ASP	A	96	26.880	65.176	13.503	1.00	11.44	7
762	CA	ASP	A	96	25.536	64.640	13.218	1.00	11.75	6
763	C	ASP	A	96	24.977	64.166	14.570	1.00	10.37	6
764	O	ASP	A	96	25.448	63.154	15.091	1.00	12.67	8
765	CB	ASP	A	96	25.622	63.467	12.217	1.00	10.69	6
766	CG	ASP	A	96	24.238	63.012	11.732	1.00	14.77	6
767	OD1	ASP	A	96	23.229	63.251	12.412	1.00	11.10	8
768	OD2	ASP	A	96	24.218	62.343	10.651	1.00	15.82	8
769	N	PHE	A	97	24.007	64.989	15.033	1.00	9.84	7
770	CA	PHE	A	97	23.482	64.615	16.377	1.00	10.91	6
771	C	PHE	A	97	22.504	63.426	16.336	1.00	13.19	6
772	O	PHE	A	97	21.945	63.125	17.373	1.00	13.15	8
773	CB	PHE	A	97	22.818	65.844	16.982	1.00	10.58	6
774	CG	PHE	A	97	23.783	66.935	17.422	1.00	14.29	6
775	CD1	PHE	A	97	24.438	67.807	16.532	1.00	13.03	6
776	CD2	PHE	A	97	23.979	67.062	18.799	1.00	13.39	6
777	CE1	PHE	A	97	25.321	68.786	17.067	1.00	11.37	6
778	CE2	PHE	A	97	24.780	68.073	19.296	1.00	9.92	6
779	CZ	PHE	A	97	25.489	68.930	18.467	1.00	9.67	6
780	N	LYS	A	98	22.220	62.889	15.152	1.00	11.59	7
781	CA	LYS	A	98	21.249	61.786	15.105	1.00	10.99	6
782	C	LYS	A	98	21.893	60.424	14.982	1.00	13.58	6
783	O	LYS	A	98	21.136	59.437	14.991	1.00	14.00	8
784	CB	LYS	A	98	20.375	61.999	13.835	1.00	10.72	6
785	CG	LYS	A	98	19.595	63.337	13.915	1.00	11.63	6
786	CD	LYS	A	98	18.627	63.408	15.091	1.00	16.21	6
787	CE	LYS	A	98	17.808	64.707	15.036	1.00	16.03	6
788	NZ	LYS	A	98	16.876	64.605	13.828	1.00	15.90	7
789	N	GLN	A	99	23.233	60.336	14.853	1.00	9.90	6
790	CA	GLN	A	99	23.900	59.079	14.613	1.00	11.53	6
791	C	GLN	A	99	25.042	58.866	15.625	1.00	12.11	6
792	O	GLN	A	99	25.610	59.842	16.118	1.00	13.41	8
793	CB	GLN	A	99	24.657	58.975	13.225	1.00	13.31	6
794	CG	GLN	A	99	23.575	58.962	12.137	1.00	20.81	6
795	CD	GLN	A	99	24.005	57.830	11.187	1.00	47.61	6
796	OE1	GLN	A	99	23.966	56.639	11.540	1.00	38.44	8
797	NE2	GLN	A	99	24.435	58.330	10.031	1.00	51.18	7
798	N	ILE	A	100	25.139	57.581	16.003	1.00	11.97	7
799	CA	ILE	A	100	26.265	57.291	16.916	1.00	9.69	6
800	C	ILE	A	100	27.586	57.395	16.160	1.00	11.29	6
801	O	ILE	A	100	27.744	57.032	14.974	1.00	13.58	8

802	CB	ILE A 100	26.080	55.834	17.385	1.00	11.20	6
803	CG1	ILE A 100	24.767	55.651	18.197	1.00	13.05	6
804	CG2	ILE A 100	27.229	55.284	18.240	1.00	10.14	6
805	CD1	ILE A 100	24.692	56.541	19.489	1.00	13.82	6
806	N	GLU A 101	28.607	57.861	16.898	1.00	11.66	7
807	CA	GLU A 101	29.968	57.886	16.322	1.00	11.19	6
808	C	GLU A 101	30.443	56.478	15.956	1.00	11.88	6
809	O	GLU A 101	30.430	55.545	16.758	1.00	12.38	8
810	CB	GLU A 101	30.918	58.478	17.400	1.00	11.53	6
811	CG	GLU A 101	32.427	58.256	17.126	1.00	10.12	6
812	CD	GLU A 101	32.796	58.791	15.715	1.00	10.82	6
813	OE1	GLU A 101	32.328	59.903	15.389	1.00	11.74	8
814	OE2	GLU A 101	33.577	58.059	15.099	1.00	12.29	8
815	N	GLU A 102	30.874	56.411	14.662	1.00	12.51	7
816	CA	GLU A 102	31.192	55.081	14.116	1.00	11.29	6
817	C	GLU A 102	32.387	54.443	14.766	1.00	12.50	6
818	O	GLU A 102	32.460	53.176	14.813	1.00	11.57	8
819	CB	GLU A 102	31.402	55.182	12.553	1.00	11.80	6
820	CG	GLU A 102	32.656	55.982	12.107	1.00	12.52	6
821	CD	GLU A 102	32.465	57.475	12.181	1.00	12.69	6
822	OE1	GLU A 102	31.368	58.018	12.431	1.00	14.09	8
823	OE2	GLU A 102	33.493	58.174	11.930	1.00	15.94	8
824	N	HIS A 103	33.391	55.145	15.324	1.00	9.73	7
825	CA	HIS A 103	34.429	54.494	16.120	1.00	9.99	6
826	C	HIS A 103	33.862	53.874	17.376	1.00	11.88	6
827	O	HIS A 103	34.531	52.943	17.864	1.00	12.86	8
828	CB	HIS A 103	35.470	55.578	16.584	1.00	12.93	6
829	CG	HIS A 103	36.364	56.049	15.481	1.00	11.28	6
830	ND1	HIS A 103	36.015	57.085	14.639	1.00	12.54	7
831	CD2	HIS A 103	37.595	55.655	15.123	1.00	14.43	6
832	CE1	HIS A 103	37.021	57.288	13.749	1.00	13.58	6
833	NE2	HIS A 103	37.965	56.434	14.054	1.00	13.95	7
834	N	PHE A 104	32.662	54.253	17.819	1.00	9.60	7
835	CA	PHE A 104	32.171	53.727	19.103	1.00	10.30	6
836	C	PHE A 104	31.116	52.625	18.924	1.00	13.76	6
837	O	PHE A 104	30.969	51.851	19.894	1.00	13.23	8
838	CB	PHE A 104	31.583	53.855	19.953	1.00	10.00	6
839	CG	PHE A 104	32.587	55.933	20.330	1.00	13.59	6
840	CD1	PHE A 104	33.954	55.782	20.204	1.00	11.92	6
841	CD2	PHE A 104	32.082	57.130	20.834	1.00	13.49	6
842	CE1	PHE A 104	34.854	56.802	20.561	1.00	12.32	6
843	CE2	PHE A 104	32.981	58.161	21.193	1.00	11.48	6
844	CZ	PHE A 104	34.365	58.005	21.053	1.00	13.46	6
845	N	GLY A 105	30.491	52.584	17.759	1.00	14.42	7
846	CA	GLY A 105	29.462	51.533	17.546	1.00	16.52	6
847	C	GLY A 105	28.362	52.143	16.670	1.00	17.63	6
848	O	GLY A 105	28.624	53.104	15.927	1.00	14.43	8
849	N	ASN A 106	27.169	51.558	16.664	1.00	14.15	7
850	CA	ASN A 106	26.017	52.150	15.960	1.00	11.27	6
851	C	ASN A 106	24.864	52.149	16.958	1.00	13.27	6
852	O	ASN A 106	25.081	51.890	18.174	1.00	13.00	8
853	CB	ASN A 106	25.756	51.332	14.677	1.00	13.55	6
854	CG	ASN A 106	25.465	49.876	14.958	1.00	18.28	6
855	OD1	ASN A 106	25.093	49.459	16.033	1.00	17.05	8
856	ND2	ASN A 106	25.576	49.021	13.910	1.00	22.96	7
857	N	TRP A 107	23.668	52.508	16.525	1.00	12.62	7
858	CA	TRP A 107	22.554	52.559	17.465	1.00	13.67	6
859	C	TRP A 107	22.296	51.203	18.121	1.00	14.10	6
860	O	TRP A 107	21.827	51.121	19.274	1.00	14.66	8
861	CB	TRP A 107	21.268	53.087	16.802	1.00	14.53	6
862	CG	TRP A 107	21.256	54.576	16.836	1.00	16.29	6
863	CD1	TRP A 107	21.351	55.357	15.696	1.00	17.56	6
864	CD2	TRP A 107	21.131	55.454	17.949	1.00	15.71	8
865	NE1	TRP A 107	21.279	56.677	16.088	1.00	16.73	7
866	CE2	TRP A 107	21.186	56.759	17.451	1.00	14.24	6
867	CE3	TRP A 107	20.995	55.256	19.345	1.00	13.30	6
868	CZ2	TRP A 107	21.082	57.915	18.240	1.00	14.74	6
869	CZ3	TRP A 107	20.919	56.410	20.131	1.00	16.80	6
870	CH2	TRP A 107	20.927	57.717	19.596	1.00	13.00	6
871	N	THR A 108	22.439	50.097	17.351	1.00	14.29	7
872	CA	THR A 108	22.316	48.792	18.020	1.00	14.34	6
873	C	THR A 108	23.305	48.617	19.154	1.00	16.18	6
874	O	THR A 108	22.945	48.095	20.224	1.00	14.57	8
875	CB	THR A 108	22.508	47.669	16.969	1.00	16.38	6
876	OG1	THR A 108	21.473	47.841	16.005	1.00	18.39	8
877	CG2	THR A 108	22.386	46.276	17.600	1.00	20.41	6
878	N	THR A 109	24.581	49.043	18.959	1.00	12.36	7
879	CA	THR A 109	25.556	48.897	20.029	1.00	13.35	6
880	C	THR A 109	25.149	49.732	21.248	1.00	12.11	6
881	O	THR A 109	25.318	49.282	22.382	1.00	12.78	8
882	CB	THR A 109	26.960	49.404	19.545	1.00	13.85	6
883	OG1	THR A 109	27.201	49.021	18.155	1.00	14.22	8
884	CG2	THR A 109	28.048	48.783	20.429	1.00	15.14	6
885	N	PHE A 110	24.673	50.945	20.982	1.00	12.96	7

#	Atom	Res	Chain	ResNum	X	Y	Z	Occ	B	El
886	CA	PHE	A	110	24.247	51.800	22.126	1.00	12.43	6
887	C	PHE	A	110	23.058	51.137	22.830	1.00	13.48	6
888	O	PHE	A	110	23.061	51.036	24.060	1.00	12.36	8
889	CB	PHE	A	110	23.823	53.160	21.525	1.00	14.28	6
890	CG	PHE	A	110	23.320	54.128	22.611	1.00	15.08	6
891	CD1	PHE	A	110	24.190	54.941	23.252	1.00	13.29	6
892	CD2	PHE	A	110	21.975	54.202	22.913	1.00	14.29	6
893	CE1	PHE	A	110	23.764	55.850	24.247	1.00	13.20	6
894	CE2	PHE	A	110	21.487	55.055	23.876	1.00	12.89	6
895	CZ	PHE	A	110	22.377	55.898	24.558	1.00	13.20	6
896	N	ASP	A	111	22.056	50.645	22.065	1.00	13.15	7
897	CA	ASP	A	111	20.916	49.993	22.755	1.00	12.63	6
898	C	ASP	A	111	21.337	48.770	23.517	1.00	14.08	6
899	O	ASP	A	111	20.917	48.606	24.698	1.00	14.41	8
900	CB	ASP	A	111	19.966	49.513	21.610	1.00	13.40	6
901	CG	ASP	A	111	19.224	50.603	20.937	1.00	18.96	6
902	OD1	ASP	A	111	19.343	51.768	21.305	1.00	17.72	8
903	OD2	ASP	A	111	18.498	50.291	19.945	1.00	18.78	8
904	N	THR	A	112	22.284	48.006	22.997	1.00	13.38	7
905	CA	THR	A	112	22.756	46.828	23.718	1.00	14.87	6
906	C	THR	A	112	23.450	47.184	25.017	1.00	14.48	6
907	O	THR	A	112	23.224	46.583	26.069	1.00	15.04	8
908	CB	THR	A	112	23.680	45.966	22.829	1.00	15.99	6
909	OG1	THR	A	112	22.844	45.644	21.711	1.00	16.67	8
910	CG2	THR	A	112	24.006	44.662	23.576	1.00	18.86	6
911	N	LEU	A	113	24.321	48.221	24.968	1.00	11.85	7
912	CA	LEU	A	113	24.982	48.631	26.219	1.00	12.44	6
913	C	LEU	A	113	23.992	49.138	27.242	1.00	12.26	6
914	O	LEU	A	113	24.057	48.793	28.403	1.00	12.86	8
915	CB	LEU	A	113	25.988	49.772	25.811	1.00	10.14	6
916	CG	LEU	A	113	26.404	50.629	27.037	1.00	13.26	6
917	CD1	LEU	A	113	27.184	49.772	28.040	1.00	13.59	6
918	CD2	LEU	A	113	27.295	51.825	26.664	1.00	12.75	6
919	N	VAL	A	114	23.020	49.971	26.823	1.00	11.82	7
920	CA	VAL	A	114	22.073	50.545	27.762	1.00	12.84	6
921	C	VAL	A	114	21.215	49.449	28.384	1.00	13.48	6
922	O	VAL	A	114	20.973	49.402	29.577	1.00	13.54	8
923	CB	VAL	A	114	21.264	51.680	27.090	1.00	13.71	6
924	CG1	VAL	A	114	20.144	52.091	28.032	1.00	18.18	6
925	CG2	VAL	A	114	22.209	52.885	26.815	1.00	15.06	6
926	N	ASN	A	115	20.760	48.534	27.512	1.00	13.50	7
927	CA	ASN	A	115	19.912	47.430	28.013	1.00	11.60	6
928	C	ASN	A	115	20.678	46.553	28.961	1.00	12.99	6
929	O	ASN	A	115	20.121	46.160	29.987	1.00	15.11	8
930	CB	ASN	A	115	19.372	46.637	26.820	1.00	15.99	6
931	CG	ASN	A	115	18.200	47.261	26.118	1.00	22.41	6
932	OD1	ASN	A	115	18.061	47.131	24.868	1.00	26.38	8
933	ND2	ASN	A	115	17.304	47.937	26.823	1.00	22.30	7
934	N	ASP	A	116	21.940	46.229	28.669	1.00	12.94	7
935	CA	ASP	A	116	22.731	45.386	29.561	1.00	12.20	6
936	C	ASP	A	116	23.087	46.144	30.835	1.00	12.69	6
937	O	ASP	A	116	23.070	45.584	31.933	1.00	12.40	8
938	CB	ASP	A	116	23.989	44.814	28.887	1.00	13.19	6
939	CG	ASP	A	116	23.648	43.698	27.896	1.00	19.33	6
940	OD1	ASP	A	116	22.461	43.487	27.582	1.00	23.80	8
941	OD2	ASP	A	116	24.583	42.967	27.460	1.00	22.90	8
942	N	ALA	A	117	23.342	47.453	30.753	1.00	12.25	7
943	CA	ALA	A	117	23.561	48.208	32.005	1.00	12.13	6
944	C	ALA	A	117	22.314	48.146	32.878	1.00	10.66	6
945	O	ALA	A	117	22.425	47.887	34.083	1.00	12.92	8
946	CB	ALA	A	117	23.877	49.692	31.625	1.00	13.05	6
947	N	HIS	A	118	21.149	48.403	32.291	1.00	12.02	7
948	CA	HIS	A	118	19.948	48.334	33.131	1.00	10.66	6
949	C	HIS	A	118	19.727	46.929	33.683	1.00	12.65	6
950	O	HIS	A	118	19.267	46.870	34.838	1.00	14.87	8
951	CB	HIS	A	118	18.714	48.622	32.200	1.00	10.79	6
952	CG	HIS	A	118	18.691	50.088	31.789	1.00	11.37	6
953	ND1	HIS	A	118	17.881	50.557	30.773	1.00	14.86	7
954	CD2	HIS	A	118	19.340	51.157	32.359	1.00	13.69	6
955	CE1	HIS	A	118	18.020	51.897	30.708	1.00	15.91	6
956	NE2	HIS	A	118	18.912	52.260	31.644	1.00	11.99	7
957	N	GLN	A	119	20.028	45.871	32.935	1.00	11.93	7
958	CA	GLN	A	119	19.843	44.545	33.592	1.00	12.22	6
959	C	GLN	A	119	20.770	44.412	34.781	1.00	15.02	6
960	O	GLN	A	119	20.519	43.615	35.694	1.00	15.33	8
961	CB	GLN	A	119	20.340	43.445	32.620	1.00	16.20	6
962	CG	GLN	A	119	19.327	43.169	31.521	1.00	17.09	6
963	CD	GLN	A	119	20.028	41.969	30.765	1.00	21.33	6
964	OE1	GLN	A	119	20.575	41.051	31.363	1.00	29.71	8
965	NE2	GLN	A	119	19.985	42.085	29.522	1.00	20.93	8
966	N	ASN	A	120	21.939	45.127	34.806	1.00	15.33	6
967	CA	ASN	A	120	22.853	45.072	35.932	1.00	16.39	6
968	C	ASN	A	120	22.541	46.148	36.970	1.00	14.25	7
969	O	ASN	A	120	23.358	46.337	37.876	1.00	15.65	8

Atom#	Atom	Res	Chain	ResNum	X	Y	Z	Occ	B	n
970	CB	ASN	A	120	24.337	45.189	35.481	1.00	12.26	6
971	CG	ASN	A	120	24.753	43.901	34.799	1.00	20.80	6
972	OD1	ASN	A	120	24.778	43.805	33.576	1.00	23.34	8
973	ND2	ASN	A	120	25.076	42.912	35.627	1.00	18.29	7
974	N	GLY	A	121	21.398	46.801	36.951	1.00	13.35	7
975	CA	GLY	A	121	20.994	47.834	37.885	1.00	16.91	6
976	C	GLY	A	121	21.840	49.129	37.772	1.00	13.99	6
977	O	GLY	A	121	21.866	49.890	38.747	1.00	15.35	8
978	N	ILE	A	122	22.262	49.397	36.527	1.00	12.67	7
979	CA	ILE	A	122	23.128	50.569	36.322	1.00	13.01	6
980	C	ILE	A	122	22.464	51.454	35.289	1.00	13.96	6
981	O	ILE	A	122	22.075	50.945	34.227	1.00	12.86	8
982	CB	ILE	A	122	24.556	50.129	35.886	1.00	12.06	6
983	CG1	ILE	A	122	25.320	49.424	37.040	1.00	15.35	6
984	CG2	ILE	A	122	25.415	51.348	35.506	1.00	13.36	6
985	CD1	ILE	A	122	26.569	48.709	36.465	1.00	15.31	6
986	N	LYS	A	123	22.344	52.752	35.609	1.00	11.45	7
987	CA	LYS	A	123	21.767	53.718	34.652	1.00	11.74	6
988	C	LYS	A	123	22.865	54.362	33.786	1.00	11.74	6
989	O	LYS	A	123	24.052	54.129	34.057	1.00	11.17	8
990	CB	LYS	A	123	21.051	54.811	35.457	1.00	11.34	6
991	CG	LYS	A	123	19.832	54.205	36.163	1.00	12.23	6
992	CD	LYS	A	123	18.994	55.310	36.815	1.00	16.30	6
993	CE	LYS	A	123	19.601	56.014	38.025	1.00	21.38	6
994	NZ	LYS	A	123	20.133	55.054	39.000	1.00	25.83	7
995	N	VAL	A	124	22.372	54.936	32.656	1.00	9.54	7
996	CA	VAL	A	124	23.343	55.533	31.740	1.00	9.53	6
997	C	VAL	A	124	22.856	56.954	31.460	1.00	12.31	6
998	O	VAL	A	124	21.723	57.168	30.990	1.00	12.23	8
999	CB	VAL	A	124	23.372	54.748	30.408	1.00	12.71	6
1000	CG1	VAL	A	124	24.327	55.480	29.398	1.00	12.93	6
1001	CG2	VAL	A	124	23.875	53.313	30.661	1.00	11.85	6
1002	N	ILE	A	125	23.726	57.937	31.756	1.00	10.85	7
1003	CA	ILE	A	125	23.419	59.311	31.352	1.00	10.22	6
1004	C	ILE	A	125	24.430	59.676	30.232	1.00	11.59	6
1005	O	ILE	A	125	25.549	59.113	30.220	1.00	12.11	8
1006	CB	ILE	A	125	23.403	60.385	32.474	1.00	10.25	6
1007	CG1	ILE	A	125	24.811	60.531	33.089	1.00	13.12	6
1008	CG2	ILE	A	125	22.304	60.035	33.484	1.00	9.99	6
1009	CD1	ILE	A	125	24.770	61.746	34.084	1.00	17.03	6
1010	N	VAL	A	126	23.971	60.450	29.252	1.00	9.62	7
1011	CA	VAL	A	126	24.864	60.725	28.102	1.00	10.33	6
1012	C	VAL	A	126	25.247	62.207	28.081	1.00	12.61	6
1013	O	VAL	A	126	24.413	63.080	28.254	1.00	11.41	8
1014	CB	VAL	A	126	24.039	60.356	26.823	1.00	11.44	6
1015	CG1	VAL	A	126	24.892	60.704	25.575	1.00	12.92	6
1016	CG2	VAL	A	126	23.704	58.883	26.856	1.00	11.52	6
1017	N	ASP	A	127	26.535	62.445	27.771	1.00	8.46	7
1018	CA	ASP	A	127	27.011	63.870	27.661	1.00	9.67	6
1019	C	ASP	A	127	26.509	64.374	26.283	1.00	10.71	6
1020	O	ASP	A	127	26.837	63.710	25.279	1.00	11.36	8
1021	CB	ASP	A	127	28.552	63.719	27.694	1.00	9.91	6
1022	CG	ASP	A	127	29.305	64.951	28.135	1.00	11.94	6
1023	OD1	ASP	A	127	28.822	66.041	27.747	1.00	11.17	8
1024	OD2	ASP	A	127	30.335	64.880	28.865	1.00	10.63	8
1025	N	PHE	A	128	25.802	65.484	26.325	1.00	8.41	7
1026	CA	PHE	A	128	25.134	65.952	25.077	1.00	9.40	6
1027	C	PHE	A	128	25.609	67.394	24.887	1.00	9.48	6
1028	O	PHE	A	128	25.752	68.163	25.858	1.00	10.67	8
1029	CB	PHE	A	128	23.609	66.006	25.437	1.00	9.59	6
1030	CG	PHE	A	128	22.760	66.405	24.223	1.00	8.93	6
1031	CD1	PHE	A	128	22.719	65.597	23.104	1.00	12.26	6
1032	CD2	PHE	A	128	22.095	67.627	24.315	1.00	10.12	6
1033	CE1	PHE	A	128	21.907	66.014	22.027	1.00	12.04	6
1034	CE2	PHE	A	128	21.325	67.993	23.183	1.00	10.05	6
1035	CZ	PHE	A	128	21.229	67.218	22.070	1.00	10.20	6
1036	N	VAL	A	129	25.985	67.736	23.647	1.00	8.13	7
1037	CA	VAL	A	129	26.851	68.932	23.398	1.00	9.78	6
1038	C	VAL	A	129	26.171	69.845	22.412	1.00	9.18	6
1039	O	VAL	A	129	26.494	70.026	21.210	1.00	10.74	8
1040	CB	VAL	A	129	28.178	68.381	22.804	1.00	11.28	6
1041	CG1	VAL	A	129	29.206	69.539	22.719	1.00	12.02	6
1042	CG2	VAL	A	129	28.829	67.246	23.590	1.00	10.62	6
1043	N	PRO	A	130	25.165	70.638	22.855	1.00	10.64	7
1044	CA	PRO	A	130	24.341	71.474	21.986	1.00	10.69	6
1045	C	PRO	A	130	24.946	72.802	21.588	1.00	11.36	6
1046	O	PRO	A	130	24.336	73.565	20.814	1.00	11.49	8
1047	CB	PRO	A	130	22.983	71.673	22.735	1.00	10.77	6
1048	CG	PRO	A	130	23.480	71.602	24.189	1.00	12.02	6
1049	CD	PRO	A	130	24.593	70.514	24.205	1.00	11.44	6
1050	N	ASN	A	131	26.107	73.181	22.144	1.00	9.48	7
1051	CA	ASN	A	131	26.687	74.481	21.859	1.00	10.43	6
1052	C	ASN	A	131	27.244	74.637	20.438	1.00	11.66	6
1053	O	ASN	A	131	27.256	75.724	19.861	1.00	11.65	8

Serial	Atom	Res	Chain	ResNum	X	Y	Z	Occ	B	
1054	CB	ASN	A	131	27.756	74.902	22.877	1.00	10.94	6
1055	CG	ASN	A	131	28.233	76.316	22.592	1.00	10.72	6
1056	OD1	ASN	A	131	27.396	77.205	22.783	1.00	10.02	8
1057	ND2	ASN	A	131	29.516	76.447	22.224	1.00	10.08	7
1058	N	HIS	A	132	27.676	73.513	19.855	1.00	10.71	7
1059	CA	HIS	A	132	28.476	73.726	18.632	1.00	9.32	6
1060	C	HIS	A	132	28.552	72.441	17.845	1.00	9.83	6
1061	O	HIS	A	132	28.256	71.362	18.361	1.00	12.46	8
1062	CB	HIS	A	132	29.896	74.227	19.005	1.00	11.52	6
1063	CG	HIS	A	132	30.560	73.394	20.080	1.00	10.70	6
1064	ND1	HIS	A	132	30.616	73.869	21.372	1.00	11.00	7
1065	CD2	HIS	A	132	31.084	72.152	20.032	1.00	9.99	6
1066	CE1	HIS	A	132	31.189	72.945	22.154	1.00	12.39	6
1067	NE2	HIS	A	132	31.445	71.929	21.368	1.00	10.00	7
1068	N	SER	A	133	28.999	72.608	16.584	1.00	10.23	7
1069	CA	SER	A	133	29.365	71.428	15.787	1.00	10.30	6
1070	C	SER	A	133	30.876	71.239	15.861	1.00	11.64	6
1071	O	SER	A	133	31.194	70.514	16.863	1.00	11.94	8
1072	CB	SER	A	133	28.807	71.514	14.344	1.00	10.65	6
1073	OG	SER	A	133	29.342	72.683	13.700	1.00	11.57	8
1074	N	THR	A	134	31.611	71.535	14.805	1.00	11.30	7
1075	CA	THR	A	134	33.034	71.082	14.740	1.00	10.51	6
1076	C	THR	A	134	33.959	72.251	14.424	1.00	8.56	6
1077	O	THR	A	134	33.553	73.376	14.121	1.00	10.40	8
1078	CB	THR	A	134	33.119	70.005	13.630	1.00	9.65	6
1079	OG1	THR	A	134	32.559	70.596	12.429	1.00	11.51	8
1080	CG2	THR	A	134	32.295	68.752	13.956	1.00	11.00	6
1081	N	PRO	A	135	35.256	71.959	14.489	1.00	10.51	7
1082	CA	PRO	A	135	36.320	72.955	14.289	1.00	12.33	6
1083	C	PRO	A	135	36.264	73.606	12.899	1.00	12.27	6
1084	O	PRO	A	135	36.014	72.968	11.868	1.00	13.19	8
1085	CB	PRO	A	135	37.627	72.145	14.405	1.00	11.11	6
1086	CG	PRO	A	135	37.241	71.129	15.486	1.00	11.56	6
1087	CD	PRO	A	135	35.809	70.746	15.111	1.00	11.98	6
1088	N	PHE	A	136	36.500	74.966	12.885	1.00	10.88	7
1089	CA	PHE	A	136	36.606	75.628	11.597	1.00	11.37	6
1090	C	PHE	A	136	37.536	76.830	11.718	1.00	11.99	6
1091	O	PHE	A	136	37.856	77.248	12.827	1.00	12.95	8
1092	CB	PHE	A	136	35.176	76.074	11.125	1.00	13.36	6
1093	CG	PHE	A	136	34.690	77.341	11.793	1.00	15.31	6
1094	CD1	PHE	A	136	34.201	77.311	13.100	1.00	12.18	6
1095	CD2	PHE	A	136	34.801	78.565	11.134	1.00	13.24	6
1096	CE1	PHE	A	136	33.813	78.477	13.738	1.00	13.44	6
1097	CE2	PHE	A	136	34.451	79.741	11.779	1.00	12.17	6
1098	CZ	PHE	A	136	33.903	79.708	13.087	1.00	15.41	6
1099	N	LYS	A	137	38.037	77.258	10.547	1.00	11.70	7
1100	CA	LYS	A	137	38.710	78.567	10.479	1.00	10.33	6
1101	C	LYS	A	137	37.859	79.484	9.618	1.00	13.10	6
1102	O	LYS	A	137	37.493	79.111	8.490	1.00	12.91	8
1103	CB	LYS	A	137	40.060	78.360	9.724	1.00	17.76	6
1104	CG	LYS	A	137	41.153	77.554	10.411	1.00	19.69	6
1105	CD	LYS	A	137	41.439	78.049	11.810	1.00	22.72	6
1106	CE	LYS	A	137	42.145	79.366	12.006	1.00	31.41	6
1107	NZ	LYS	A	137	43.212	79.771	11.035	1.00	25.39	7
1108	N	ALA	A	138	37.657	80.688	10.158	1.00	12.09	7
1109	CA	ALA	A	138	36.683	81.510	9.375	1.00	12.57	6
1110	C	ALA	A	138	37.267	81.873	8.017	1.00	14.60	6
1111	O	ALA	A	138	36.469	82.176	7.094	1.00	14.01	8
1112	CB	ALA	A	138	36.410	82.806	10.148	1.00	15.45	6
1113	N	ASN	A	139	38.597	81.986	7.900	1.00	13.25	7
1114	CA	ASN	A	139	39.165	82.359	6.608	1.00	14.41	6
1115	C	ASN	A	139	39.444	81.212	5.682	1.00	14.25	6
1116	O	ASN	A	139	40.047	81.349	4.562	1.00	14.82	8
1117	CB	ASN	A	139	40.443	83.188	6.852	1.00	16.78	6
1118	CG	ASN	A	139	41.666	82.292	7.083	1.00	22.27	6
1119	OD1	ASN	A	139	41.484	81.167	7.486	1.00	25.66	8
1120	ND2	ASN	A	139	42.853	82.762	6.773	1.00	21.53	7
1121	N	ASP	A	140	39.150	79.955	6.052	1.00	14.76	7
1122	CA	ASP	A	140	39.433	78.835	5.203	1.00	13.60	6
1123	C	ASP	A	140	38.470	77.674	5.337	1.00	16.09	6
1124	O	ASP	A	140	38.607	76.905	6.339	1.00	13.95	8
1125	CB	ASP	A	140	40.885	78.382	5.557	1.00	12.14	6
1126	CG	ASP	A	140	41.331	77.260	4.643	1.00	16.87	6
1127	OD1	ASP	A	140	40.616	76.705	3.810	1.00	17.82	8
1128	OD2	ASP	A	140	42.569	76.954	4.814	1.00	25.09	8
1129	N	SER	A	141	37.529	77.555	4.414	1.00	15.21	7
1130	CA	SER	A	141	36.508	76.520	4.501	1.00	16.63	6
1131	C	SER	A	141	37.048	75.092	4.285	1.00	17.44	6
1132	O	SER	A	141	36.349	74.129	4.607	1.00	18.28	8
1133	CB	SER	A	141	35.372	76.746	3.493	1.00	19.52	6
1134	OG	SER	A	141	35.867	76.579	2.144	1.00	16.38	8
1135	N	THR	A	142	38.302	74.958	3.839	1.00	14.08	7
1136	CA	THR	A	142	38.889	73.615	3.649	1.00	15.63	6
1137	C	THR	A	142	39.445	73.036	4.933	1.00	16.91	6

Serial	Atom	Res	Chain	ResNum	X	Y	Z	Occ	B	Elem
1138	O	THR	A	142	39.798	71.840	4.961	1.00	15.67	8
1139	CB	THR	A	142	40.105	73.658	2.622	1.00	16.69	6
1140	OG1	THR	A	142	41.261	74.288	3.211	1.00	18.35	8
1141	CG2	THR	A	142	39.653	74.352	1.368	1.00	25.03	6
1142	N	PHE	A	143	39.533	73.874	6.007	1.00	15.48	7
1143	CA	PHE	A	143	40.027	73.314	7.285	1.00	12.99	6
1144	C	PHE	A	143	38.953	72.384	7.908	1.00	14.57	6
1145	O	PHE	A	143	37.807	72.837	8.026	1.00	12.49	8
1146	CB	PHE	A	143	40.412	74.481	8.238	1.00	14.72	6
1147	CG	PHE	A	143	40.905	73.939	9.546	1.00	13.33	6
1148	CD1	PHE	A	143	42.192	73.454	9.669	1.00	14.66	6
1149	CD2	PHE	A	143	40.054	73.908	10.679	1.00	10.60	6
1150	CE1	PHE	A	143	42.677	72.946	10.878	1.00	14.87	6
1151	CE2	PHE	A	143	40.556	73.381	11.849	1.00	12.09	6
1152	CZ	PHE	A	143	41.842	72.912	11.975	1.00	16.30	6
1153	N	ALA	A	144	39.342	71.174	8.264	1.00	14.83	7
1154	CA	ALA	A	144	38.380	70.250	8.884	1.00	14.49	6
1155	C	ALA	A	144	37.165	70.136	7.976	1.00	14.94	6
1156	O	ALA	A	144	37.369	69.878	6.784	1.00	13.97	8
1157	CB	ALA	A	144	37.990	70.683	10.323	1.00	12.28	6
1158	N	GLU	A	145	35.942	70.135	8.506	1.00	11.43	7
1159	CA	GLU	A	145	34.744	70.061	7.645	1.00	10.19	6
1160	C	GLU	A	145	34.063	71.386	7.520	1.00	10.66	6
1161	O	GLU	A	145	32.824	71.513	7.266	1.00	11.84	8
1162	CB	GLU	A	145	33.771	68.943	8.180	1.00	11.91	6
1163	CG	GLU	A	145	34.408	67.577	8.042	1.00	11.68	6
1164	CD	GLU	A	145	33.591	66.467	8.697	1.00	15.77	6
1165	OE1	GLU	A	145	32.530	66.660	9.208	1.00	20.95	8
1166	OE2	GLU	A	145	34.122	65.351	8.743	1.00	21.22	8
1167	N	GLY	A	146	34.677	72.533	7.898	1.00	12.55	7
1168	CA	GLY	A	146	34.021	73.826	7.783	1.00	12.03	6
1169	C	GLY	A	146	32.799	73.976	8.739	1.00	15.18	6
1170	O	GLY	A	146	32.025	74.880	8.511	1.00	13.95	8
1171	N	GLY	A	147	32.774	73.157	9.790	1.00	12.57	7
1172	CA	GLY	A	147	31.639	73.265	10.703	1.00	12.28	6
1173	C	GLY	A	147	30.439	72.415	10.267	1.00	14.09	6
1174	O	GLY	A	147	29.372	72.644	10.903	1.00	11.87	8
1175	N	ALA	A	148	30.552	71.583	9.258	1.00	11.40	7
1176	CA	ALA	A	148	29.343	70.942	8.707	1.00	13.48	6
1177	C	ALA	A	148	28.495	70.132	9.729	1.00	11.58	6
1178	O	ALA	A	148	29.025	69.390	10.557	1.00	12.14	8
1179	CB	ALA	A	148	29.861	69.927	7.660	1.00	14.52	6
1180	N	LEU	A	149	27.188	70.305	9.543	1.00	12.19	7
1181	CA	LEU	A	149	26.192	69.594	10.365	1.00	10.39	6
1182	C	LEU	A	149	25.298	68.798	9.407	1.00	10.91	6
1183	O	LEU	A	149	24.907	69.311	8.313	1.00	14.52	8
1184	CB	LEU	A	149	25.345	70.684	11.028	1.00	10.56	6
1185	CG	LEU	A	149	24.344	70.222	12.098	1.00	14.34	6
1186	CD1	LEU	A	149	25.067	69.618	13.294	1.00	16.21	6
1187	CD2	LEU	A	149	23.436	71.409	12.447	1.00	15.60	6
1188	N	TYR	A	150	25.047	67.560	9.836	1.00	12.69	7
1189	CA	TYR	A	150	24.178	66.675	9.057	1.00	11.91	6
1190	C	TYR	A	150	23.007	66.234	9.957	1.00	13.54	6
1191	O	TYR	A	150	23.116	66.203	11.188	1.00	12.53	8
1192	CB	TYR	A	150	24.945	65.392	8.671	1.00	12.62	6
1193	CG	TYR	A	150	26.104	65.699	7.694	1.00	10.80	6
1194	CD1	TYR	A	150	27.275	66.240	8.185	1.00	13.28	6
1195	CD2	TYR	A	150	25.983	65.330	6.359	1.00	14.10	6
1196	CE1	TYR	A	150	28.295	66.545	7.273	1.00	14.81	6
1197	CE2	TYR	A	150	27.004	65.609	5.479	1.00	17.28	6
1198	CZ	TYR	A	150	28.139	66.206	5.959	1.00	17.89	8
1199	OH	TYR	A	150	29.227	66.444	5.082	1.00	18.31	6
1200	N	ASN	A	151	21.932	65.787	5.288	1.00	11.59	7
1201	CA	ASN	A	151	20.774	65.261	10.048	1.00	13.13	6
1202	C	ASN	A	151	20.582	63.801	9.659	1.00	14.46	6
1203	O	ASN	A	151	20.020	63.527	8.595	1.00	15.41	8
1204	CB	ASN	A	151	19.542	66.068	9.633	1.00	12.78	6
1205	CG	ASN	A	151	18.280	65.607	10.386	1.00	15.85	6
1206	OD1	ASN	A	151	18.376	65.084	11.460	1.00	14.91	8
1207	ND2	ASN	A	151	17.115	65.851	9.790	1.00	23.13	7
1208	N	ASN	A	152	21.153	62.881	10.455	1.00	11.88	7
1209	CA	ASN	A	152	21.160	61.460	10.056	1.00	12.91	6
1210	C	ASN	A	152	21.628	61.265	8.619	1.00	17.10	6
1211	O	ASN	A	152	21.059	60.495	7.804	1.00	18.15	8
1212	CB	ASN	A	152	19.763	60.894	10.305	1.00	15.18	6
1213	CG	ASN	A	152	19.772	59.363	10.289	1.00	28.25	6
1214	OD1	ASN	A	152	20.803	58.741	10.579	1.00	25.62	8
1215	ND2	ASN	A	152	18.647	58.722	9.925	1.00	26.09	7
1216	N	GLY	A	153	22.797	61.857	8.344	1.00	11.86	6
1217	CA	GLY	A	153	23.494	61.698	7.061	1.00	15.63	6
1218	C	GLY	A	153	23.096	62.727	6.007	1.00	17.70	6
1219	O	GLY	A	153	23.819	62.836	4.994	1.00	16.41	8
1220	N	THR	A	154	21.975	63.414	6.209	1.00	13.77	7
1221	CA	THR	A	154	21.535	64.406	5.220	1.00	14.28	6

#	Atom	Res	Chain	Seq	X	Y	Z	Occ	B	El
1222	C	THR	A	154	22.181	65.781	5.538	1.00	14.24	6
1223	O	THR	A	154	22.048	66.264	6.650	1.00	15.31	8
1224	CB	THR	A	154	20.008	64.592	5.212	1.00	21.10	6
1225	OG1	THR	A	154	19.488	63.334	4.709	1.00	20.66	8
1226	CG2	THR	A	154	19.569	65.711	4.260	1.00	21.99	6
1227	N	TYR	A	155	22.977	66.280	4.566	1.00	14.57	7
1228	CA	TYR	A	155	23.613	67.573	4.928	1.00	15.65	6
1229	C	TYR	A	155	22.652	68.698	5.184	1.00	17.55	6
1230	O	TYR	A	155	21.639	68.912	4.487	1.00	16.44	8
1231	CB	TYR	A	155	24.440	67.984	3.678	1.00	16.08	6
1232	CG	TYR	A	155	25.238	69.243	3.820	1.00	16.71	6
1233	CD1	TYR	A	155	26.324	69.277	4.693	1.00	16.20	6
1234	CD2	TYR	A	155	24.989	70.381	3.075	1.00	16.54	6
1235	CE1	TYR	A	155	27.139	70.407	4.789	1.00	16.44	6
1236	CE2	TYR	A	155	25.773	71.530	3.162	1.00	14.17	6
1237	CZ	TYR	A	155	26.843	71.515	4.015	1.00	16.83	6
1238	OH	TYR	A	155	27.673	72.601	4.142	1.00	15.80	8
1239	N	MET	A	156	22.895	69.456	6.272	1.00	12.36	7
1240	CA	MET	A	156	22.120	70.658	6.595	1.00	12.72	6
1241	C	MET	A	156	22.877	71.939	6.202	1.00	14.70	6
1242	O	MET	A	156	22.290	72.876	5.628	1.00	14.86	8
1243	CB	MET	A	156	21.886	70.683	8.141	1.00	14.72	6
1244	CG	MET	A	156	21.045	69.510	8.543	1.00	13.16	6
1245	SD	MET	A	156	20.812	69.391	10.354	1.00	16.44	16
1246	CE	MET	A	156	19.828	70.788	10.735	1.00	16.14	6
1247	N	GLY	A	157	24.138	72.003	6.634	1.00	15.28	7
1248	CA	GLY	A	157	24.888	73.257	6.260	1.00	12.22	6
1249	C	GLY	A	157	26.169	73.337	7.061	1.00	15.29	6
1250	O	GLY	A	157	26.402	72.513	7.965	1.00	13.96	8
1251	N	ASN	A	158	26.981	74.369	6.736	1.00	13.19	7
1252	CA	ASN	A	158	28.205	74.586	7.485	1.00	12.06	6
1253	C	ASN	A	158	28.353	76.085	7.759	1.00	8.80	6
1254	O	ASN	A	158	27.377	76.850	7.583	1.00	11.73	8
1255	CB	ASN	A	158	29.438	73.957	6.787	1.00	12.79	6
1256	CG	ASN	A	158	29.783	74.647	5.457	1.00	16.08	6
1257	OD1	ASN	A	158	29.311	75.727	5.160	1.00	12.69	8
1258	ND2	ASN	A	158	30.650	74.060	4.603	1.00	20.88	7
1259	N	TYR	A	159	29.559	75.484	8.260	1.00	10.67	7
1260	CA	TYR	A	159	29.714	77.924	8.640	1.00	11.67	6
1261	C	TYR	A	159	29.665	78.827	7.432	1.00	12.87	6
1262	O	TYR	A	159	29.444	80.029	7.605	1.00	12.70	8
1263	CB	TYR	A	159	31.055	78.072	9.434	1.00	11.86	6
1264	CG	TYR	A	159	31.171	79.475	10.063	1.00	11.61	6
1265	CD1	TYR	A	159	30.414	79.868	11.149	1.00	14.86	6
1266	CD2	TYR	A	159	31.962	80.402	9.380	1.00	13.87	6
1267	CE1	TYR	A	159	30.576	81.181	11.600	1.00	15.19	6
1268	CE2	TYR	A	159	32.069	81.716	9.811	1.00	19.88	6
1269	CZ	TYR	A	159	31.387	82.088	10.926	1.00	16.15	6
1270	OH	TYR	A	159	31.377	83.383	11.432	1.00	16.24	8
1271	N	PHE	A	160	29.904	78.307	6.253	1.00	13.52	7
1272	CA	PHE	A	160	30.073	79.126	5.029	1.00	12.84	6
1273	C	PHE	A	160	28.855	79.182	4.153	1.00	12.34	6
1274	O	PHE	A	160	28.803	80.101	3.300	1.00	14.63	8
1275	CB	PHE	A	160	31.228	78.421	4.234	1.00	12.99	6
1276	CG	PHE	A	160	32.504	78.508	5.080	1.00	14.21	6
1277	CD1	PHE	A	160	33.310	79.662	4.935	1.00	12.64	6
1278	CD2	PHE	A	160	32.804	77.488	5.965	1.00	12.84	6
1279	CE1	PHE	A	160	34.466	79.772	5.737	1.00	14.91	6
1280	CE2	PHE	A	160	33.940	77.621	6.761	1.00	14.27	6
1281	CZ	PHE	A	160	34.769	78.740	6.653	1.00	13.93	6
1282	N	ASP	A	161	27.917	78.220	4.232	1.00	14.16	7
1283	CA	ASP	A	161	26.731	78.329	3.371	1.00	14.46	6
1284	C	ASP	A	161	25.486	78.622	4.217	1.00	14.25	6
1285	O	ASP	A	161	24.375	78.215	3.808	1.00	15.23	8
1286	CB	ASP	A	161	26.557	77.031	2.565	1.00	12.66	6
1287	CG	ASP	A	161	26.500	75.766	3.373	1.00	15.72	6
1288	OD1	ASP	A	161	26.191	75.896	4.579	1.00	13.04	8
1289	OD2	ASP	A	161	26.767	74.716	2.766	1.00	20.26	8
1290	N	ASP	A	162	25.656	79.460	5.227	1.00	13.84	7
1291	CA	ASP	A	162	24.557	79.667	6.196	1.00	14.60	6
1292	C	ASP	A	162	23.787	80.939	5.979	1.00	16.54	6
1293	O	ASP	A	162	22.840	81.252	6.726	1.00	18.25	8
1294	CB	ASP	A	162	25.253	79.785	7.584	1.00	12.11	6
1295	CG	ASP	A	162	24.264	79.524	8.717	1.00	13.12	6
1296	OD1	ASP	A	162	23.408	78.624	8.551	1.00	12.76	8
1297	OD2	ASP	A	162	24.417	80.201	9.775	1.00	12.76	8
1298	N	ALA	A	163	24.109	81.725	4.949	1.00	16.09	7
1299	CA	ALA	A	163	23.428	83.019	4.792	1.00	16.11	6
1300	C	ALA	A	163	21.914	82.901	4.848	1.00	17.25	6
1301	O	ALA	A	163	21.341	83.825	5.467	1.00	24.30	8
1302	CB	ALA	A	163	23.828	83.609	3.414	1.00	19.57	6
1303	N	THR	A	164	21.317	81.987	4.145	1.00	19.25	7
1304	CA	THR	A	164	19.845	81.973	4.121	1.00	24.11	6
1305	C	THR	A	164	19.237	81.014	5.149	1.00	25.86	6

1306	O	THR A 164	18.055	80.605	5.002	1.00	25.38	8
1307	CB	THR A 164	19.384	81.502	2.723	1.00	21.33	6
1308	OG1	THR A 164	19.834	80.146	2.496	1.00	29.22	8
1309	CG2	THR A 164	20.062	82.359	1.658	1.00	32.67	6
1310	N	LYS A 165	20.086	80.431	6.008	1.00	19.16	7
1311	CA	LYS A 165	19.577	79.430	6.929	1.00	16.62	6
1312	C	LYS A 165	19.714	79.888	8.391	1.00	18.03	6
1313	O	LYS A 165	18.735	79.767	9.173	1.00	17.25	8
1314	CB	LYS A 165	20.423	78.140	6.826	1.00	13.80	6
1315	CG	LYS A 165	20.215	77.497	5.423	1.00	21.49	6
1316	CD	LYS A 165	20.913	76.186	5.334	1.00	30.47	6
1317	CE	LYS A 165	22.394	76.297	5.245	1.00	26.09	6
1318	NZ	LYS A 165	23.067	75.308	4.349	1.00	18.44	7
1319	N	GLY A 166	20.839	80.499	8.700	1.00	14.27	7
1320	CA	GLY A 166	21.048	81.024	10.082	1.00	14.43	6
1321	C	GLY A 166	21.103	79.863	11.109	1.00	14.80	6
1322	O	GLY A 166	20.730	80.096	12.287	1.00	14.92	8
1323	N	TYR A 167	21.732	78.765	10.734	1.00	12.78	7
1324	CA	TYR A 167	21.882	77.692	11.732	1.00	12.54	6
1325	C	TYR A 167	22.991	78.009	12.739	1.00	12.62	6
1326	O	TYR A 167	23.085	77.349	13.776	1.00	11.78	8
1327	CB	TYR A 167	22.226	76.374	11.022	1.00	10.92	6
1328	CG	TYR A 167	21.127	75.836	10.103	1.00	14.63	6
1329	CD1	TYR A 167	19.810	76.232	10.291	1.00	13.59	6
1330	CD2	TYR A 167	21.490	74.921	9.121	1.00	15.06	6
1331	CE1	TYR A 167	18.836	75.722	9.421	1.00	13.53	6
1332	CE2	TYR A 167	20.503	74.385	8.254	1.00	12.93	6
1333	CZ	TYR A 167	19.211	74.809	8.469	1.00	15.10	6
1334	OH	TYR A 167	18.235	74.289	7.594	1.00	18.52	8
1335	N	PHE A 168	23.963	78.870	12.342	1.00	11.36	7
1336	CA	PHE A 168	25.072	79.208	13.244	1.00	11.20	6
1337	C	PHE A 168	25.097	80.677	13.551	1.00	12.02	6
1338	O	PHE A 168	24.515	81.539	12.854	1.00	11.21	8
1339	CB	PHE A 168	26.432	78.934	12.493	1.00	13.41	6
1340	CG	PHE A 168	26.552	77.459	12.174	1.00	11.52	6
1341	CD1	PHE A 168	27.044	76.583	13.130	1.00	10.63	6
1342	CD2	PHE A 168	26.171	77.007	10.899	1.00	13.81	6
1343	CE1	PHE A 168	27.122	75.214	12.765	1.00	12.04	6
1344	CE2	PHE A 168	26.250	75.639	10.574	1.00	12.05	6
1345	CZ	PHE A 168	26.752	74.751	11.518	1.00	12.13	6
1346	N	HIS A 169	25.665	81.067	14.709	1.00	11.43	7
1347	CA	HIS A 169	25.979	82.473	14.979	1.00	12.52	6
1348	C	HIS A 169	27.266	82.836	14.195	1.00	11.69	6
1349	O	HIS A 169	28.109	82.035	13.921	1.00	11.66	8
1350	CB	HIS A 169	26.361	82.658	16.482	1.00	11.57	6
1351	CG	HIS A 169	25.157	82.347	17.376	1.00	11.97	6
1352	ND1	HIS A 169	25.403	81.657	18.588	1.00	11.36	7
1353	CD2	HIS A 169	23.838	82.604	17.274	1.00	11.55	6
1354	CE1	HIS A 169	24.195	81.518	19.195	1.00	12.02	6
1355	NE2	HIS A 169	23.233	82.111	18.427	1.00	10.80	8
1356	N	HIS A 170	27.295	84.147	13.797	1.00	11.77	7
1357	CA	HIS A 170	28.474	84.629	13.015	1.00	12.20	6
1358	C	HIS A 170	29.029	85.872	13.672	1.00	14.93	6
1359	O	HIS A 170	29.174	86.974	13.053	1.00	16.92	8
1360	CB	HIS A 170	28.083	84.949	11.533	1.00	13.28	6
1361	CG	HIS A 170	27.535	83.698	10.888	1.00	12.02	6
1362	ND1	HIS A 170	28.327	82.925	10.069	1.00	15.51	7
1363	CD2	HIS A 170	26.306	83.088	10.915	1.00	13.12	6
1364	CE1	HIS A 170	27.639	81.863	9.689	1.00	16.69	6
1365	NE2	HIS A 170	26.409	81.953	10.156	1.00	13.32	7
1366	N	ASN A 171	29.387	85.778	14.962	1.00	12.58	7
1367	CA	ASN A 171	29.735	86.967	15.733	1.00	12.76	6
1368	C	ASN A 171	31.201	87.040	16.147	1.00	13.34	6
1369	O	ASN A 171	31.554	87.947	16.949	1.00	17.31	8
1370	CB	ASN A 171	28.916	86.961	17.054	1.00	13.40	6
1371	CG	ASN A 171	27.430	86.948	16.719	1.00	17.12	6
1372	OD1	ASN A 171	26.595	86.168	17.252	1.00	16.54	8
1373	ND2	ASN A 171	27.046	87.866	15.861	1.00	12.82	7
1374	N	GLY A 172	32.013	86.197	15.601	1.00	13.14	7
1375	CA	GLY A 172	33.444	86.166	15.964	1.00	15.42	6
1376	C	GLY A 172	33.728	85.275	17.210	1.00	16.69	6
1377	O	GLY A 172	32.817	84.706	17.722	1.00	14.98	8
1378	N	ASP A 173	34.993	85.180	17.526	1.00	14.76	7
1379	CA	ASP A 173	35.473	84.346	18.622	1.00	12.93	6
1380	C	ASP A 173	35.292	84.996	19.976	1.00	12.76	6
1381	O	ASP A 173	35.410	86.248	20.131	1.00	11.79	8
1382	CB	ASP A 173	36.980	84.152	18.369	1.00	14.28	6
1383	CG	ASP A 173	37.273	83.139	17.268	1.00	24.98	6
1384	OD1	ASP A 173	36.398	82.387	16.822	1.00	17.11	8
1385	OD2	ASP A 173	38.451	83.124	16.815	1.00	23.71	8
1386	N	ILE A 174	35.073	84.127	20.969	1.00	12.58	7
1387	CA	ILE A 174	35.136	84.670	22.362	1.00	11.65	6
1388	C	ILE A 174	36.500	85.307	22.646	1.00	14.87	6
1389	O	ILE A 174	37.508	84.670	22.337	1.00	15.09	8

#	Atom	Res	Ch	Seq	X	Y	Z	Occ	B	E
1390	CB	ILE	A	174	34.896	83.495	23.357	1.00	13.27	6
1391	CG1	ILE	A	174	33.431	83.005	23.177	1.00	10.95	6
1392	CG2	ILE	A	174	35.145	84.016	24.806	1.00	12.74	6
1393	CD1	ILE	A	174	33.220	81.690	24.000	1.00	11.77	6
1394	N	SER	A	175	36.441	86.493	23.260	1.00	14.22	7
1395	CA	SER	A	175	37.710	87.093	23.770	1.00	16.89	6
1396	C	SER	A	175	37.712	87.131	25.291	1.00	18.02	6
1397	O	SER	A	175	38.617	86.587	25.938	1.00	19.20	8
1398	CB	SER	A	175	37.868	88.470	23.138	1.00	17.86	6
1399	OG	SER	A	175	39.049	89.044	23.724	1.00	24.28	8
1400	N	ASN	A	176	36.650	87.662	25.854	1.00	14.39	7
1401	CA	ASN	A	176	36.515	87.678	27.336	1.00	13.41	6
1402	C	ASN	A	176	35.511	86.561	27.678	1.00	12.76	6
1403	O	ASN	A	176	34.286	86.760	27.482	1.00	13.43	8
1404	CB	ASN	A	176	35.898	89.032	27.724	1.00	15.61	6
1405	CG	ASN	A	176	35.749	89.123	29.243	1.00	17.91	6
1406	OD1	ASN	A	176	35.963	88.166	29.982	1.00	15.18	8
1407	ND2	ASN	A	176	35.402	90.347	29.694	1.00	22.13	7
1408	N	TRP	A	177	36.085	85.465	28.237	1.00	14.29	7
1409	CA	TRP	A	177	35.172	84.361	28.558	1.00	13.39	6
1410	C	TRP	A	177	34.248	84.677	29.724	1.00	15.05	6
1411	O	TRP	A	177	33.279	83.898	29.909	1.00	14.25	8
1412	CB	TRP	A	177	36.054	83.145	28.953	1.00	16.12	6
1413	CG	TRP	A	177	36.712	82.559	27.721	1.00	14.43	6
1414	CD1	TRP	A	177	37.745	83.101	26.998	1.00	16.58	6
1415	CD2	TRP	A	177	36.399	81.291	27.142	1.00	13.92	6
1416	NE1	TRP	A	177	38.070	82.235	25.940	1.00	18.57	7
1417	CE2	TRP	A	177	37.234	81.131	26.014	1.00	18.68	6
1418	CE3	TRP	A	177	35.437	80.298	27.392	1.00	17.34	6
1419	CZ2	TRP	A	177	37.148	80.031	25.169	1.00	14.28	6
1420	CZ3	TRP	A	177	35.379	79.182	26.574	1.00	17.20	6
1421	CH2	TRP	A	177	36.253	79.045	25.441	1.00	18.00	6
1422	N	ASP	A	178	34.477	85.795	30.469	1.00	12.43	7
1423	CA	ASP	A	178	33.507	86.126	31.517	1.00	11.08	6
1424	C	ASP	A	178	32.454	87.115	31.053	1.00	12.88	6
1425	O	ASP	A	178	31.586	87.420	31.881	1.00	15.73	8
1426	CB	ASP	A	178	34.243	86.717	32.739	1.00	17.78	6
1427	CG	ASP	A	178	35.201	85.739	33.362	1.00	24.33	6
1428	OD1	ASP	A	178	34.916	84.535	33.440	1.00	18.07	8
1429	OD2	ASP	A	178	36.317	86.155	33.777	1.00	24.77	8
1430	N	ASP	A	179	32.527	87.608	29.810	1.00	11.79	7
1431	CA	ASP	A	179	31.448	88.502	29.357	1.00	11.42	6
1432	C	ASP	A	179	30.292	87.603	28.903	1.00	12.33	6
1433	O	ASP	A	179	30.537	86.725	28.067	1.00	12.08	8
1434	CB	ASP	A	179	31.997	89.341	28.184	1.00	15.43	6
1435	CG	ASP	A	179	30.831	90.225	27.763	1.00	18.80	6
1436	OD1	ASP	A	179	30.462	91.232	28.431	1.00	20.56	8
1437	OD2	ASP	A	179	30.213	89.893	26.753	1.00	13.51	8
1438	N	ARG	A	180	29.115	87.768	29.477	1.00	13.02	7
1439	CA	ARG	A	180	28.057	86.771	29.205	1.00	11.54	6
1440	C	ARG	A	180	27.585	86.833	27.757	1.00	11.36	6
1441	O	ARG	A	180	27.261	85.746	27.225	1.00	10.36	8
1442	CB	ARG	A	180	26.893	87.051	30.172	1.00	11.27	6
1443	CG	ARG	A	180	27.286	86.796	31.654	1.00	13.31	6
1444	CD	ARG	A	180	27.797	85.340	31.899	1.00	12.49	6
1445	NE	ARG	A	180	26.694	84.389	31.571	1.00	10.75	8
1446	CZ	ARG	A	180	26.896	83.305	30.812	1.00	11.38	6
1447	NH1	ARG	A	180	28.090	82.893	30.359	1.00	11.55	7
1448	NH2	ARG	A	180	25.769	82.597	30.589	1.00	11.21	8
1449	N	TYR	A	181	27.508	87.990	27.113	1.00	13.64	7
1450	CA	TYR	A	181	27.104	87.980	25.688	1.00	10.51	6
1451	C	TYR	A	181	28.195	87.277	24.870	1.00	10.68	6
1452	O	TYR	A	181	27.826	86.403	24.044	1.00	11.14	8
1453	CB	TYR	A	181	26.915	89.446	25.196	1.00	10.66	6
1454	CG	TYR	A	181	26.645	89.417	23.698	1.00	12.30	6
1455	CD1	TYR	A	181	25.446	89.009	23.179	1.00	12.16	6
1456	CD2	TYR	A	181	27.712	89.736	22.837	1.00	13.07	6
1457	CE1	TYR	A	181	25.242	88.936	21.808	1.00	15.09	6
1458	CE2	TYR	A	181	27.510	89.688	21.457	1.00	18.03	6
1459	CZ	TYR	A	181	26.275	89.265	20.988	1.00	16.90	7
1460	OH	TYR	A	181	26.097	89.156	19.614	1.00	19.76	8
1461	N	GLU	A	182	29.473	87.528	25.083	1.00	17.30	7
1462	CA	GLU	A	182	30.468	86.836	24.265	1.00	13.50	6
1463	C	GLU	A	182	30.442	85.311	24.504	1.00	12.35	6
1464	O	GLU	A	182	30.482	84.533	23.582	1.00	10.45	8
1465	CB	GLU	A	182	31.939	87.266	24.571	1.00	11.45	6
1466	CG	GLU	A	182	32.131	88.769	24.214	1.00	10.48	6
1467	CD	GLU	A	182	33.640	89.046	24.246	1.00	12.66	8
1468	OE1	GLU	A	182	34.487	88.229	23.958	1.00	19.30	8
1469	OE2	GLU	A	182	34.009	90.199	24.690	1.00	15.48	8
1470	N	ALA	A	183	30.314	84.945	25.796	1.00	10.95	7
1471	CA	ALA	A	183	30.436	83.521	26.110	1.00	10.17	6
1472	C	ALA	A	183	29.302	82.709	25.471	1.00	11.02	6
1473	O	ALA	A	183	29.555	81.542	25.197	1.00	10.91	8

#	Atom	Res	Ch	Seq	X	Y	Z	Occ	B	E
1474	CB	ALA	A	183	30.290	83.352	27.664	1.00	11.34	6
1475	N	GLN	A	184	28.196	83.381	25.172	1.00	10.70	7
1476	CA	GLN	A	184	27.046	82.624	24.601	1.00	8.62	6
1477	C	GLN	A	184	26.939	82.844	23.102	1.00	10.69	6
1478	O	GLN	A	184	26.509	81.913	22.388	1.00	10.72	8
1479	CB	GLN	A	184	25.772	83.078	25.330	1.00	12.35	6
1480	CG	GLN	A	184	25.730	82.584	26.785	1.00	9.80	6
1481	CD	GLN	A	184	24.603	83.315	27.538	1.00	12.56	6
1482	OE1	GLN	A	184	24.739	84.527	27.890	1.00	15.21	8
1483	NE2	GLN	A	184	23.536	82.580	27.775	1.00	8.78	7
1484	N	TRP	A	185	27.186	84.018	22.585	1.00	11.90	7
1485	CA	TRP	A	185	26.968	84.285	21.148	1.00	9.74	6
1486	C	TRP	A	185	28.252	84.298	20.318	1.00	10.30	6
1487	O	TRP	A	185	28.093	84.277	19.065	1.00	11.67	8
1488	CB	TRP	A	185	26.201	85.647	20.965	1.00	12.61	6
1489	CG	TRP	A	105	24.696	85.390	21.039	1.00	10.95	6
1490	CD1	TRP	A	185	23.863	85.166	19.989	1.00	12.59	6
1491	CD2	TRP	A	185	23.898	85.345	22.226	1.00	12.07	6
1492	NE1	TRP	A	185	22.561	84.887	20.428	1.00	14.15	7
1493	CE2	TRP	A	185	22.600	85.003	21.805	1.00	13.63	6
1494	CE3	TRP	A	185	24.154	84.846	23.587	1.00	13.34	6
1495	CZ2	TRP	A	185	21.534	85.361	22.703	1.00	14.01	6
1496	CZ3	TRP	A	185	23.083	85.004	24.494	1.00	14.75	6
1497	CH2	TRP	A	185	21.812	84.254	24.035	1.00	13.91	6
1498	N	LYS	A	186	29.413	84.170	20.924	1.00	11.31	7
1499	CA	LYS	A	186	30.655	82.764	20.127	1.00	12.29	6
1500	C	LYS	A	186	31.228	81.896	20.238	1.00	14.72	6
1501	O	LYS	A	186	30.718	82.003	20.981	1.00	12.47	8
1502	CB	LYS	A	186	31.682	85.224	20.582	1.00	11.09	6
1503	CG	LYS	A	186	31.145	86.645	20.243	1.00	14.74	6
1504	CD	LYS	A	186	32.295	87.655	20.601	1.00	14.53	6
1505	CE	LYS	A	186	31.688	89.051	20.286	1.00	20.29	6
1506	NZ	LYS	A	186	32.744	89.943	19.726	1.00	28.67	7
1507	N	ASN	A	187	32.217	82.446	19.379	1.00	13.16	7
1508	CA	ASN	A	187	32.653	81.074	19.195	1.00	13.22	6
1509	C	ASN	A	187	33.587	80.604	20.314	1.00	15.43	6
1510	O	ASN	A	187	34.587	81.250	20.646	1.00	12.76	8
1511	CB	ASN	A	187	33.386	80.850	17.862	1.00	12.72	6
1512	CG	ASN	A	187	32.673	81.537	16.697	1.00	18.84	6
1513	OD1	ASN	A	187	31.447	81.456	16.631	1.00	15.46	8
1514	ND2	ASN	A	187	33.426	82.219	15.839	1.00	16.94	7
1515	N	PHE	A	188	33.369	79.356	20.719	1.00	11.91	7
1516	CA	PHE	A	188	34.321	78.658	21.604	1.00	10.93	6
1517	C	PHE	A	188	35.600	78.457	20.806	1.00	13.95	6
1518	O	PHE	A	188	35.534	78.128	19.632	1.00	13.55	8
1519	CB	PHE	A	188	33.682	77.259	21.822	1.00	11.05	6
1520	CG	PHE	A	188	34.177	76.552	23.071	1.00	11.08	6
1521	CD1	PHE	A	188	35.431	75.953	23.060	1.00	14.39	6
1522	CD2	PHE	A	188	33.364	76.492	24.194	1.00	14.52	6
1523	CE1	PHE	A	188	35.865	75.309	24.204	1.00	15.63	6
1524	CE2	PHE	A	188	33.821	75.818	25.327	1.00	14.28	6
1525	CZ	PHE	A	188	35.081	75.233	25.350	1.00	12.75	6
1526	N	THR	A	189	36.737	78.710	21.504	1.00	11.89	7
1527	CA	THR	A	189	38.000	78.578	20.769	1.00	11.77	6
1528	C	THR	A	189	38.851	77.524	21.457	1.00	13.65	6
1529	O	THR	A	189	38.630	77.149	22.589	1.00	14.67	8
1530	CB	THR	A	189	38.826	79.904	20.788	1.00	12.56	6
1531	OG1	THR	A	189	39.066	80.215	22.180	1.00	15.52	8
1532	CG2	THR	A	189	38.012	81.045	20.136	1.00	13.79	6
1533	N	ASP	A	190	39.773	76.961	20.639	1.00	11.10	7
1534	CA	ASP	A	190	40.736	75.985	21.186	1.00	11.48	6
1535	C	ASP	A	190	42.109	76.575	20.929	1.00	11.63	6
1536	O	ASP	A	190	42.403	77.103	19.861	1.00	13.03	8
1537	CB	ASP	A	190	40.530	74.703	20.365	1.00	11.07	6
1538	CG	ASP	A	190	41.445	73.591	20.781	1.00	12.55	6
1539	OD1	ASP	A	190	42.691	73.716	20.943	1.00	14.12	8
1540	OD2	ASP	A	190	40.937	72.422	20.956	1.00	14.32	8
1541	N	PRO	A	191	43.013	76.539	21.885	1.00	12.39	7
1542	CA	PRO	A	191	44.344	77.094	21.756	1.00	16.15	6
1543	C	PRO	A	191	45.205	76.488	20.648	1.00	17.03	6
1544	O	PRO	A	191	46.194	77.139	20.258	1.00	17.33	8
1545	CB	PRO	A	191	45.077	76.806	23.067	1.00	15.82	6
1546	CG	PRO	A	191	43.951	76.489	24.024	1.00	19.99	6
1547	CD	PRO	A	191	42.769	75.960	23.220	1.00	15.14	6
1548	N	ALA	A	192	44.778	75.374	20.102	1.00	12.41	7
1549	CA	ALA	A	192	45.446	74.851	18.871	1.00	16.29	6
1550	C	ALA	A	192	45.279	75.807	17.697	1.00	21.11	6
1551	O	ALA	A	192	46.014	75.779	16.675	1.00	19.85	8
1552	CB	ALA	A	192	44.978	73.466	18.579	1.00	20.48	6
1553	N	GLY	A	193	44.317	76.695	17.671	1.00	16.93	7
1554	CA	GLY	A	193	44.199	77.733	16.641	1.00	17.85	6
1555	C	GLY	A	193	42.919	77.582	15.819	1.00	17.03	6
1556	O	GLY	A	193	42.991	77.960	14.651	1.00	17.03	8
1557	N	PHE	A	194	41.888	76.955	16.373	1.00	13.47	7

Atom#	Atom	Res	Chain	Seq	X	Y	Z	Occ	B	El
1558	CA	PHE	A	194	40.612	76.976	15.567	1.00	11.84	6
1559	C	PHE	A	194	39.441	77.265	16.536	1.00	11.03	6
1560	O	PHE	A	194	39.621	77.423	17.750	1.00	11.69	8
1561	CB	PHE	A	194	40.411	75.629	14.855	1.00	10.98	6
1562	CG	PHE	A	194	40.568	74.412	15.767	1.00	11.46	6
1563	CD1	PHE	A	194	39.545	74.063	16.649	1.00	11.56	6
1564	CD2	PHE	A	194	41.707	73.656	15.747	1.00	15.24	6
1565	CE1	PHE	A	194	39.688	72.942	17.460	1.00	11.26	6
1566	CE2	PHE	A	194	41.871	72.533	16.574	1.00	12.49	6
1567	CZ	PHE	A	194	40.860	72.181	17.450	1.00	12.03	6
1568	N	SER	A	195	38.283	77.497	15.895	1.00	10.73	7
1569	CA	SER	A	195	37.097	77.782	16.704	1.00	11.52	6
1570	C	SER	A	195	36.081	76.649	16.423	1.00	11.92	6
1571	O	SER	A	195	36.284	75.902	15.519	1.00	10.36	8
1572	CB	SER	A	195	36.416	79.059	16.186	1.00	16.30	6
1573	OG	SER	A	195	37.442	80.119	16.216	1.00	22.79	8
1574	N	LEU	A	196	35.060	75.594	17.262	1.00	12.22	7
1575	CA	LEU	A	196	34.007	76.384	17.060	1.00	12.16	6
1576	C	LEU	A	196	32.756	77.416	16.584	1.00	7.60	6
1577	O	LEU	A	196	32.364	74.883	17.112	1.00	13.16	8
1578	CB	LEU	A	196	33.660	74.298	18.410	1.00	10.95	6
1579	CG	LEU	A	196	34.880	73.454	19.107	1.00	10.75	6
1580	CD1	LEU	A	196	34.439	73.384	20.334	1.00	11.59	6
1581	CD2	LEU	A	196	35.719	72.300	18.134	1.00	12.48	6
1582	N	ALA	A	197	32.139	75.877	15.481	1.00	9.97	7
1583	CA	ALA	A	197	30.995	76.637	14.905	1.00	10.64	6
1584	C	ALA	A	197	29.788	76.658	15.830	1.00	14.16	6
1585	O	ALA	A	197	29.362	75.622	16.314	1.00	11.22	8
1586	CB	ALA	A	197	30.629	75.928	13.565	1.00	10.95	6
1587	N	ASP	A	198	29.429	77.869	16.236	1.00	10.28	7
1588	CA	ASP	A	198	28.459	78.009	17.350	1.00	10.50	6
1589	C	ASP	A	198	27.030	77.880	16.795	1.00	11.99	6
1590	O	ASP	A	198	26.607	78.731	15.993	1.00	13.33	8
1591	CB	ASP	A	198	28.744	79.433	17.900	1.00	12.29	6
1592	CG	ASP	A	198	28.236	79.529	19.353	1.00	10.72	6
1593	OD1	ASP	A	198	28.683	78.683	20.172	1.00	11.33	8
1594	OD2	ASP	A	198	27.401	80.452	19.671	1.00	10.45	8
1595	N	LEU	A	199	26.310	76.847	17.247	1.00	10.29	7
1596	CA	LEU	A	199	24.927	76.690	16.763	1.00	8.80	6
1597	C	LEU	A	199	24.086	77.821	17.321	1.00	9.88	6
1598	O	LEU	A	199	24.246	78.320	18.449	1.00	11.27	8
1599	CB	LEU	A	199	24.452	75.296	17.279	1.00	9.52	6
1600	CG	LEU	A	199	25.178	74.135	16.526	1.00	10.88	6
1601	CD1	LEU	A	199	24.799	72.879	17.321	1.00	12.59	6
1602	CD2	LEU	A	199	24.658	74.045	15.071	1.00	14.19	6
1603	N	SER	A	200	23.024	78.144	16.505	1.00	11.57	7
1604	CA	SER	A	200	22.055	79.160	16.950	1.00	10.26	6
1605	C	SER	A	200	20.810	78.424	17.499	1.00	11.50	6
1606	O	SER	A	200	19.994	77.917	16.741	1.00	11.79	8
1607	CB	SER	A	200	21.636	80.012	15.731	1.00	14.77	6
1608	OG	SER	A	200	20.723	81.011	16.249	1.00	13.69	8
1609	N	GLN	A	201	20.786	78.294	18.837	1.00	10.61	7
1610	CA	GLN	A	201	19.599	77.668	19.473	1.00	11.55	6
1611	C	GLN	A	201	18.421	78.648	19.371	1.00	11.79	6
1612	O	GLN	A	201	17.305	78.161	19.700	1.00	12.43	8
1613	CB	GLN	A	201	19.852	77.359	20.969	1.00	11.94	6
1614	CG	GLN	A	201	21.042	76.370	21.151	1.00	10.05	6
1615	CD	GLN	A	201	22.393	77.086	21.126	1.00	10.84	6
1616	OE1	GLN	A	201	22.499	78.298	21.208	1.00	11.60	8
1617	NE2	GLN	A	201	23.465	76.231	21.079	1.00	9.83	7
1618	N	GLU	A	202	18.590	79.860	18.862	1.00	10.41	7
1619	CA	GLU	A	202	17.434	80.763	18.667	1.00	11.55	6
1620	C	GLU	A	202	16.849	80.531	17.277	1.00	14.51	6
1621	O	GLU	A	202	15.856	81.166	16.903	1.00	16.57	8
1622	CB	GLU	A	202	17.877	82.226	18.875	1.00	10.69	6
1623	CG	GLU	A	202	18.522	82.442	20.252	1.00	9.01	6
1624	CD	GLU	A	202	20.002	82.069	20.330	1.00	12.78	6
1625	OE1	GLU	A	202	20.680	82.031	19.299	1.00	16.62	8
1626	OE2	GLU	A	202	20.457	81.782	21.434	1.00	12.31	8
1627	N	ASN	A	203	17.507	79.704	16.435	1.00	11.04	7
1628	CA	ASN	A	203	16.939	79.318	15.142	1.00	10.64	6
1629	C	ASN	A	203	16.020	78.098	15.372	1.00	13.48	6
1630	O	ASN	A	203	16.441	77.148	16.008	1.00	12.81	8
1631	CB	ASN	A	203	18.105	78.892	14.217	1.00	11.71	6
1632	CG	ASN	A	203	17.604	78.307	12.930	1.00	15.36	6
1633	OD1	ASN	A	203	17.271	77.124	12.838	1.00	15.37	8
1634	ND2	ASN	A	203	17.611	79.108	11.829	1.00	15.33	7
1635	N	GLY	A	204	14.797	78.219	14.831	1.00	13.94	7
1636	CA	GLY	A	204	13.813	77.115	15.175	1.00	14.96	6
1637	C	GLY	A	204	14.202	75.766	14.593	1.00	12.98	6
1638	O	GLY	A	204	13.891	74.775	15.243	1.00	14.33	8
1639	N	THR	A	205	14.802	75.705	13.401	1.00	11.88	7
1640	CA	THR	A	205	15.279	74.405	12.894	1.00	12.45	6
1641	C	THR	A	205	16.275	73.780	13.856	1.00	11.35	6

#	Atom	Res	Chain	Seq	X	Y	Z	Occ	B	n
1642	O	THR	A	205	16.161	72.621	14.172	1.00	11.84	8
1643	CB	THR	A	205	15.866	74.603	11.497	1.00	14.40	6
1644	OG1	THR	A	205	14.760	75.060	10.662	1.00	18.35	8
1645	CG2	THR	A	205	16.344	73.256	10.930	1.00	16.31	6
1646	N	ILE	A	206	17.295	74.586	14.188	1.00	10.89	7
1647	CA	ILE	A	206	18.330	73.996	15.081	1.00	10.54	6
1648	C	ILE	A	206	17.736	73.689	16.457	1.00	9.48	6
1649	O	ILE	A	206	18.081	72.638	17.051	1.00	10.61	8
1650	CB	ILE	A	206	19.481	75.017	15.212	1.00	9.16	6
1651	CG1	ILE	A	206	20.193	75.155	13.844	1.00	12.29	6
1652	CG2	ILE	A	206	20.550	74.624	16.273	1.00	11.91	6
1653	CD1	ILE	A	206	20.691	73.847	13.218	1.00	12.69	6
1654	N	ALA	A	207	16.912	74.562	17.021	1.00	10.30	7
1655	CA	ALA	A	207	16.350	74.185	18.339	1.00	12.45	6
1656	C	ALA	A	207	15.583	72.871	18.311	1.00	12.96	6
1657	O	ALA	A	207	15.714	72.053	19.217	1.00	13.37	8
1658	CB	ALA	A	207	15.511	75.359	18.883	1.00	13.30	6
1659	N	GLN	A	208	14.749	72.676	17.282	1.00	10.55	7
1660	CA	GLN	A	208	13.968	71.431	17.236	1.00	11.60	6
1661	C	GLN	A	208	14.842	70.256	16.877	1.00	14.22	6
1662	O	GLN	A	208	14.627	69.200	17.426	1.00	11.79	8
1663	CB	GLN	A	208	12.869	71.573	16.136	1.00	13.62	6
1664	CG	GLN	A	208	11.847	70.429	16.197	1.00	14.37	6
1665	CD	GLN	A	208	11.089	70.392	17.513	1.00	15.09	6
1666	OE1	GLN	A	208	10.565	71.371	17.957	1.00	14.67	8
1667	NE2	GLN	A	208	11.168	69.230	18.180	1.00	14.23	7
1668	N	TYR	A	209	15.876	70.454	16.062	1.00	11.41	7
1669	CA	TYR	A	209	16.807	69.372	15.724	1.00	10.71	6
1670	C	TYR	A	209	17.570	68.930	16.979	1.00	10.31	6
1671	O	TYR	A	209	17.634	67.727	17.231	1.00	11.12	8
1672	CB	TYR	A	209	17.840	69.989	14.743	1.00	11.54	6
1673	CG	TYR	A	209	19.072	69.148	14.457	1.00	12.56	6
1674	CD1	TYR	A	209	19.031	68.083	13.584	1.00	11.38	6
1675	CD2	TYR	A	209	20.269	69.485	15.060	1.00	12.87	6
1676	CE1	TYR	A	209	20.184	67.341	13.286	1.00	16.27	8
1677	CE2	TYR	A	209	21.432	68.739	14.794	1.00	13.55	6
1678	CZ	TYR	A	209	21.368	67.676	13.915	1.00	13.50	6
1679	OH	TYR	A	209	22.527	66.968	13.654	1.00	14.21	6
1680	N	LEU	A	210	17.985	69.883	17.821	1.00	10.88	7
1681	CA	LEU	A	210	18.731	69.461	19.027	1.00	10.76	6
1682	C	LEU	A	210	17.776	68.829	20.049	1.00	11.52	6
1683	O	LEU	A	210	18.178	67.863	20.687	1.00	11.55	8
1684	CB	LEU	A	210	19.388	70.688	19.713	1.00	10.40	6
1685	CG	LEU	A	210	20.603	71.237	18.893	1.00	10.20	6
1686	CD1	LEU	A	210	20.918	72.628	19.480	1.00	12.47	6
1687	CD2	LEU	A	210	21.766	70.260	18.920	1.00	13.21	6
1688	N	THR	A	211	16.539	69.327	20.069	1.00	9.66	7
1689	CA	THR	A	211	15.520	68.689	20.933	1.00	11.98	6
1690	C	THR	A	211	15.294	67.272	20.496	1.00	11.93	6
1691	O	THR	A	211	15.311	66.354	21.335	1.00	11.42	8
1692	CB	THR	A	211	14.193	69.502	20.889	1.00	12.29	6
1693	OG1	THR	A	211	14.488	70.842	21.373	1.00	11.73	8
1694	CG2	THR	A	211	13.137	68.832	21.808	1.00	10.37	6
1695	N	ASP	A	212	15.071	67.084	19.170	1.00	12.47	7
1696	CA	ASP	A	212	14.813	65.738	18.671	1.00	10.73	6
1697	C	ASP	A	212	15.998	64.812	18.997	1.00	12.42	6
1698	O	ASP	A	212	15.748	63.598	19.194	1.00	12.22	8
1699	CB	ASP	A	212	14.605	65.829	17.159	1.00	12.33	6
1700	CG	ASP	A	212	13.253	66.477	16.797	1.00	15.18	6
1701	OD1	ASP	A	212	12.379	66.643	17.667	1.00	13.61	8
1702	OD2	ASP	A	212	13.100	66.828	15.585	1.00	14.64	8
1703	N	ALA	A	213	17.230	65.326	18.830	1.00	10.02	7
1704	CA	ALA	A	213	18.376	64.422	19.084	1.00	10.72	6
1705	C	ALA	A	213	18.422	64.023	20.552	1.00	10.18	6
1706	O	ALA	A	213	18.819	62.882	20.861	1.00	11.96	8
1707	CB	ALA	A	213	19.679	65.163	18.735	1.00	10.83	6
1708	N	ALA	A	214	18.155	64.970	21.448	1.00	12.55	7
1709	CA	ALA	A	214	18.185	64.599	22.892	1.00	10.81	6
1710	C	ALA	A	214	17.073	63.631	23.235	1.00	11.35	6
1711	O	ALA	A	214	17.246	62.667	23.953	1.00	12.75	8
1712	CB	ALA	A	214	18.038	65.876	23.757	1.00	10.21	6
1713	N	VAL	A	215	15.885	63.854	22.677	1.00	10.82	7
1714	CA	VAL	A	215	14.724	62.923	22.875	1.00	11.60	6
1715	C	VAL	A	215	15.035	61.577	22.302	1.00	13.68	6
1716	O	VAL	A	215	14.673	60.552	22.903	1.00	15.03	8
1717	CB	VAL	A	215	13.462	63.523	22.283	1.00	14.89	6
1718	CG1	VAL	A	215	12.285	62.514	22.234	1.00	16.68	6
1719	CG2	VAL	A	215	12.982	64.740	23.099	1.00	15.78	6
1720	N	GLN	A	216	15.759	61.496	21.193	1.00	12.25	7
1721	CA	GLN	A	216	16.153	60.192	20.632	1.00	13.52	6
1722	C	GLN	A	216	17.023	59.404	21.577	1.00	13.62	6
1723	O	GLN	A	216	16.864	58.192	21.706	1.00	12.75	8
1724	CB	GLN	A	216	16.814	60.382	19.232	1.00	12.82	6
1725	CG	GLN	A	216	17.225	59.032	18.639	1.00	14.36	6

#	Atom	Residue	Chain	Seq	X	Y	Z	Occ	B	n
1726	CD	GLN	A	216	17.856	59.162	17.261	1.00	18.32	6
1727	OE1	GLN	A	216	18.762	59.976	16.980	1.00	21.98	8
1728	NE2	GLN	A	216	17.392	58.348	16.335	1.00	19.07	7
1729	N	LEU	A	217	17.996	60.043	22.269	1.00	11.16	7
1730	CA	LEU	A	217	18.781	59.317	23.261	1.00	11.02	6
1731	C	LEU	A	217	17.885	58.756	24.396	1.00	10.47	6
1732	O	LEU	A	217	18.106	57.631	24.754	1.00	13.19	8
1733	CB	LEU	A	217	19.847	60.252	23.876	1.00	12.00	6
1734	CG	LEU	A	217	20.974	60.577	22.859	1.00	10.07	6
1735	CD1	LEU	A	217	21.789	61.789	23.418	1.00	10.56	6
1736	CD2	LEU	A	217	21.953	59.383	22.787	1.00	12.18	6
1737	N	VAL	A	218	16.940	59.584	24.844	1.00	10.50	7
1738	CA	VAL	A	218	16.027	59.030	25.884	1.00	14.12	6
1739	C	VAL	A	218	15.114	57.918	25.304	1.00	12.90	6
1740	O	VAL	A	218	14.914	56.898	25.978	1.00	14.29	8
1741	CB	VAL	A	218	15.121	60.168	26.376	1.00	11.91	6
1742	CG1	VAL	A	218	14.131	59.671	27.428	1.00	14.89	6
1743	CG2	VAL	A	218	16.045	61.169	27.107	1.00	14.59	6
1744	N	ALA	A	219	14.717	57.992	24.051	1.00	14.93	7
1745	CA	ALA	A	219	13.868	56.915	23.478	1.00	14.21	6
1746	C	ALA	A	219	14.647	55.619	23.377	1.00	16.64	6
1747	O	ALA	A	219	14.072	54.517	23.401	1.00	14.94	8
1748	CB	ALA	A	219	13.379	57.270	22.059	1.00	15.54	6
1749	N	HIS	A	220	15.959	55.702	23.258	1.00	12.54	7
1750	CA	HIS	A	220	16.853	54.570	23.215	1.00	14.41	6
1751	C	HIS	A	220	17.305	54.124	24.611	1.00	12.25	6
1752	O	HIS	A	220	18.194	53.243	24.711	1.00	13.99	8
1753	CB	HIS	A	220	18.055	54.802	22.293	1.00	11.69	6
1754	CG	HIS	A	220	17.630	54.760	20.840	1.00	14.69	6
1755	ND1	HIS	A	220	17.984	53.694	20.064	1.00	14.71	7
1756	CD2	HIS	A	220	16.928	55.624	20.077	1.00	16.42	6
1757	CE1	HIS	A	220	17.518	53.916	18.823	1.00	16.42	6
1758	NE2	HIS	A	220	16.855	55.077	18.794	1.00	17.85	7
1759	N	GLY	A	221	16.735	54.666	25.655	1.00	11.47	7
1760	CA	GLY	A	221	16.964	54.129	26.999	1.00	11.21	6
1761	C	GLY	A	221	17.655	55.060	27.965	1.00	13.22	6
1762	O	GLY	A	221	17.875	54.635	29.095	1.00	13.24	8
1763	N	ALA	A	222	18.297	56.111	27.426	1.00	12.29	7
1764	CA	ALA	A	222	19.139	56.912	28.356	1.00	14.03	6
1765	C	ALA	A	222	18.325	57.381	29.553	1.00	13.32	6
1766	O	ALA	A	222	17.138	57.794	29.459	1.00	10.61	8
1767	CB	ALA	A	222	19.700	58.136	27.618	1.00	12.95	6
1768	N	ASP	A	223	19.009	57.376	30.709	1.00	11.67	7
1769	CA	ASP	A	223	18.391	57.768	31.951	1.00	10.83	6
1770	C	ASP	A	223	18.621	59.249	32.306	1.00	12.58	6
1771	O	ASP	A	223	18.116	59.667	33.356	1.00	11.47	8
1772	CB	ASP	A	223	18.997	56.964	33.123	1.00	9.67	6
1773	CG	ASP	A	223	18.744	55.469	32.925	1.00	12.72	6
1774	OD1	ASP	A	223	17.554	55.095	33.224	1.00	15.35	8
1775	OD2	ASP	A	223	19.610	54.686	32.482	1.00	10.77	8
1776	N	GLY	A	224	19.186	59.972	31.348	1.00	11.00	7
1777	CA	GLY	A	224	19.393	61.413	31.651	1.00	11.62	6
1778	C	GLY	A	224	20.641	61.861	30.809	1.00	9.62	6
1779	O	GLY	A	224	21.069	61.107	29.928	1.00	9.69	8
1780	N	LEU	A	225	20.869	63.123	30.983	1.00	10.32	7
1781	CA	LEU	A	225	21.962	63.749	30.177	1.00	9.63	6
1782	C	LEU	A	225	22.828	64.566	31.128	1.00	10.99	6
1783	O	LEU	A	225	22.356	65.110	32.097	1.00	10.86	8
1784	CB	LEU	A	225	21.389	64.780	29.172	1.00	9.89	6
1785	CG	LEU	A	225	20.424	64.212	28.122	1.00	10.00	6
1786	CD1	LEU	A	225	19.806	65.361	27.279	1.00	12.69	6
1787	CD2	LEU	A	225	21.003	63.092	27.243	1.00	13.97	6
1788	N	ARG	A	226	24.135	64.671	30.790	1.00	10.52	7
1789	CA	ARG	A	226	24.993	65.777	31.295	1.00	9.01	6
1790	C	ARG	A	226	25.093	66.744	30.110	1.00	10.97	6
1791	O	ARG	A	226	25.628	66.348	29.083	1.00	10.94	8
1792	CB	ARG	A	226	26.337	65.159	31.691	1.00	9.10	6
1793	CG	ARG	A	226	27.381	66.213	32.158	1.00	8.52	6
1794	CD	ARG	A	226	28.248	66.648	30.956	1.00	9.82	6
1795	NE	ARG	A	226	29.438	67.400	31.425	1.00	9.22	7
1796	CZ	ARG	A	226	30.251	68.074	30.592	1.00	9.27	6
1797	NH1	ARG	A	226	29.978	68.191	29.289	1.00	9.46	7
1798	NH2	ARG	A	226	31.311	68.687	31.114	1.00	10.94	7
1799	N	ILE	A	227	24.590	67.940	30.262	1.00	10.46	7
1800	CA	ILE	A	227	24.553	68.901	29.124	1.00	7.87	6
1801	C	ILE	A	227	25.807	69.751	29.238	1.00	8.86	6
1802	O	ILE	A	227	26.042	70.450	30.199	1.00	11.29	8
1803	CB	ILE	A	227	23.295	69.783	29.269	1.00	9.62	6
1804	CG1	ILE	A	227	22.096	68.814	29.378	1.00	10.28	6
1805	CG2	ILE	A	227	23.196	70.601	27.964	1.00	11.21	6
1806	CD1	ILE	A	227	20.743	69.547	29.412	1.00	15.31	6
1807	N	ASP	A	228	26.544	69.672	28.125	1.00	9.30	7
1808	CA	ASP	A	228	27.846	70.399	28.079	1.00	9.30	6
1809	C	ASP	A	228	27.672	71.915	27.904	1.00	9.50	6

1810	O	ASP	A	228	26.757	72.338	27.189	1.00	9.88	8
1811	CB	ASP	A	228	28.521	69.865	26.769	1.00	9.23	6
1812	CG	ASP	A	228	29.904	70.442	26.587	1.00	10.24	6
1813	OD1	ASP	A	228	30.725	70.228	27.505	1.00	10.28	8
1814	OD2	ASP	A	228	30.208	71.128	25.574	1.00	9.89	8
1815	N	ALA	A	229	28.575	72.633	28.582	1.00	8.93	7
1816	CA	ALA	A	229	28.745	74.068	28.257	1.00	9.72	6
1817	C	ALA	A	229	27.455	74.868	28.324	1.00	10.58	6
1818	O	ALA	A	229	27.123	75.695	27.464	1.00	10.73	8
1819	CB	ALA	A	229	29.355	74.180	26.841	1.00	9.90	6
1820	N	VAL	A	230	26.729	74.708	29.487	1.00	9.05	7
1821	CA	VAL	A	230	25.421	75.392	29.548	1.00	9.09	6
1822	C	VAL	A	230	25.547	76.905	29.753	1.00	10.83	6
1823	O	VAL	A	230	24.587	77.636	29.573	1.00	12.48	8
1824	CB	VAL	A	230	24.469	74.836	30.634	1.00	10.80	6
1825	CG1	VAL	A	230	24.119	73.389	30.262	1.00	9.84	6
1826	CG2	VAL	A	230	25.084	74.934	32.047	1.00	11.19	6
1827	N	LYS	A	231	26.753	77.312	30.189	1.00	9.10	7
1828	CA	LYS	A	231	26.988	78.763	30.234	1.00	7.91	6
1829	C	LYS	A	231	27.304	79.391	28.883	1.00	11.01	6
1830	O	LYS	A	231	27.398	80.604	28.809	1.00	13.02	8
1831	CB	LYS	A	231	28.173	79.043	31.209	1.00	10.48	6
1832	CG	LYS	A	231	29.567	78.801	30.602	1.00	10.63	6
1833	CD	LYS	A	231	30.599	78.817	31.758	1.00	13.75	6
1834	CE	LYS	A	231	30.937	80.205	32.236	1.00	13.70	6
1835	NZ	LYS	A	231	32.215	80.157	33.104	1.00	11.49	7
1836	N	HIS	A	232	27.364	78.589	27.814	1.00	9.37	7
1837	CA	HIS	A	232	27.744	79.081	26.480	1.00	12.72	6
1838	C	HIS	A	232	26.576	79.044	25.537	1.00	12.48	6
1839	O	HIS	A	232	26.698	79.505	24.388	1.00	9.62	8
1840	CB	HIS	A	232	28.849	78.129	25.915	1.00	9.94	6
1841	CG	HIS	A	232	30.145	78.345	26.709	1.00	9.35	6
1842	ND1	HIS	A	232	30.780	79.598	26.674	1.00	11.03	7
1843	CD2	HIS	A	232	30.803	77.530	27.554	1.00	11.85	6
1844	CE1	HIS	A	232	31.838	79.483	27.513	1.00	12.22	6
1845	NE2	HIS	A	232	31.882	78.243	28.049	1.00	11.68	7
1846	N	PHE	A	233	25.342	78.757	25.941	1.00	9.67	7
1847	CA	PHE	A	233	24.165	78.957	25.062	1.00	10.57	6
1848	C	PHE	A	233	23.015	79.437	25.991	1.00	9.21	6
1849	O	PHE	A	233	23.073	79.254	27.225	1.00	11.50	8
1850	CB	PHE	A	233	23.792	77.705	24.296	1.00	9.50	6
1851	CG	PHE	A	233	23.382	76.523	25.145	1.00	10.66	6
1852	CD1	PHE	A	233	24.335	75.588	25.559	1.00	12.13	6
1853	CD2	PHE	A	233	22.023	76.331	25.466	1.00	11.40	6
1854	CE1	PHE	A	233	23.982	74.503	26.347	1.00	12.18	6
1855	CE2	PHE	A	233	21.642	75.219	26.237	1.00	10.50	6
1856	CZ	PHE	A	233	22.629	74.340	26.657	1.00	11.82	6
1857	N	ASN	A	234	22.051	80.148	25.349	1.00	12.13	7
1858	CA	ASN	A	234	21.093	80.799	26.253	1.00	8.31	6
1859	C	ASN	A	234	20.367	79.859	27.214	1.00	9.45	6
1860	O	ASN	A	234	20.112	78.722	26.829	1.00	9.94	8
1861	CB	ASN	A	234	20.132	81.662	25.369	1.00	10.71	6
1862	CG	ASN	A	234	18.981	80.871	24.740	1.00	12.93	6
1863	OD1	ASN	A	234	18.070	80.516	25.519	1.00	12.50	8
1864	ND2	ASN	A	234	18.975	80.590	23.448	1.00	10.79	7
1865	N	SER	A	235	20.023	80.451	28.374	1.00	10.78	7
1866	CA	SER	A	235	19.401	79.586	29.396	1.00	8.67	6
1867	C	SER	A	235	17.906	79.375	29.177	1.00	11.06	6
1868	O	SER	A	235	17.399	78.402	29.759	1.00	11.56	8
1869	CB	ASER	A	235	19.594	80.196	30.792	0.60	11.41	6
1870	OG	ASER	A	235	20.974	80.269	31.098	0.60	10.81	8
1869	CB	BSER	A	235	19.679	80.067	30.817	0.40	10.03	6
1870	OG	BSER	A	235	19.311	81.423	30.914	0.40	7.85	8
1871	N	GLY	A	236	17.301	80.140	28.312	1.00	10.81	7
1872	CA	GLY	A	236	15.906	79.857	27.882	1.00	11.83	6
1873	C	GLY	A	236	15.821	78.448	27.289	1.00	12.34	6
1874	O	GLY	A	236	14.942	77.638	27.552	1.00	11.07	8
1875	N	PHE	A	237	16.742	78.158	26.370	1.00	10.75	7
1876	CA	PHE	A	237	16.793	76.864	25.702	1.00	9.60	6
1877	C	PHE	A	237	17.034	75.738	26.715	1.00	10.90	6
1878	O	PHE	A	237	16.476	74.641	26.526	1.00	10.74	8
1879	CB	PHE	A	237	17.714	76.891	24.467	1.00	12.14	6
1880	CG	PHE	A	237	17.642	75.595	23.691	1.00	12.02	6
1881	CD1	PHE	A	237	16.442	75.242	23.048	1.00	12.55	6
1882	CD2	PHE	A	237	18.751	74.740	23.659	1.00	11.64	6
1883	CE1	PHE	A	237	16.367	74.046	22.358	1.00	12.60	6
1884	CE2	PHE	A	237	18.634	73.533	22.952	1.00	11.39	6
1885	CZ	PHE	A	237	17.468	73.200	22.301	1.00	13.43	6
1886	N	SER	A	238	17.965	75.939	27.658	1.00	10.56	7
1887	CA	SER	A	238	18.119	74.844	28.637	1.00	11.43	6
1888	C	SER	A	238	16.762	74.489	29.289	1.00	10.24	6
1889	O	SER	A	238	16.416	73.312	29.442	1.00	10.58	8
1890	CB	SER	A	238	19.069	75.303	29.776	1.00	11.86	6
1891	OG	SER	A	238	20.432	75.070	29.404	1.00	12.16	6

1892	N	LYS	A	239	16.053	75.523	29.757	1.00	10.06	7
1893	CA	LYS	A	239	14.749	75.262	30.433	1.00	10.43	6
1894	C	LYS	A	239	13.712	74.684	29.505	1.00	10.95	6
1895	O	LYS	A	239	13.006	73.718	29.879	1.00	11.91	8
1896	CB	LYS	A	239	14.259	76.598	31.123	1.00	8.56	6
1897	CG	LYS	A	239	12.889	76.329	31.861	1.00	12.08	6
1898	CD	LYS	A	239	12.577	77.644	32.648	1.00	11.14	6
1899	CE	LYS	A	239	11.131	77.442	33.240	1.00	11.14	6
1900	NZ	LYS	A	239	10.797	78.668	34.098	1.00	10.40	7
1901	N	SER	A	240	13.600	75.153	28.279	1.00	9.61	7
1902	CA	SER	A	240	12.611	74.583	27.330	1.00	9.08	6
1903	C	SER	A	240	13.006	73.207	26.957	1.00	10.90	6
1904	O	SER	A	240	12.160	72.320	26.790	1.00	10.77	8
1905	CB	SER	A	240	12.560	75.572	26.136	1.00	11.38	6
1906	OG	SER	A	240	11.488	75.039	25.266	1.00	12.68	8
1907	N	LEU	A	241	14.300	72.896	26.747	1.00	10.43	7
1908	CA	LEU	A	241	14.726	71.560	26.389	1.00	10.29	6
1909	C	LEU	A	241	14.420	70.599	27.539	1.00	11.73	6
1910	O	LEU	A	241	13.924	69.507	27.303	1.00	11.15	8
1911	CB	LEU	A	241	16.255	71.582	26.077	1.00	10.07	6
1912	CG	LEU	A	241	16.816	70.170	25.829	1.00	12.71	6
1913	CD1	LEU	A	241	16.205	69.502	24.577	1.00	13.39	6
1914	CD2	LEU	A	241	18.333	70.350	25.611	1.00	12.90	6
1915	N	ALA	A	242	14.710	71.055	28.778	1.00	11.25	7
1916	CA	ALA	A	242	14.427	70.181	29.912	1.00	10.06	6
1917	C	ALA	A	242	12.923	69.852	29.969	1.00	10.68	6
1918	O	ALA	A	242	12.565	68.703	30.215	1.00	10.97	8
1919	CB	ALA	A	242	14.910	70.889	31.196	1.00	10.49	6
1920	N	ASP	A	243	12.062	70.870	29.712	1.00	10.40	7
1921	CA	ASP	A	243	10.609	70.595	29.694	1.00	11.54	6
1922	C	ASP	A	243	10.365	69.448	28.700	1.00	12.78	6
1923	O	ASP	A	243	9.636	68.474	29.006	1.00	12.31	8
1924	CB	ASP	A	243	9.930	71.904	29.186	1.00	11.95	6
1925	CG	ASP	A	243	8.507	71.717	28.674	1.00	13.21	6
1926	OD1	ASP	A	243	7.668	71.113	29.422	1.00	12.98	8
1927	OD2	ASP	A	243	8.223	72.178	27.546	1.00	12.92	8
1928	N	LYS	A	244	10.825	69.536	27.451	1.00	10.42	7
1929	CA	LYS	A	244	10.523	68.449	26.484	1.00	10.94	6
1930	C	LYS	A	244	10.997	67.097	26.998	1.00	13.31	6
1931	O	LYS	A	244	10.349	66.061	26.763	1.00	12.24	8
1932	CB	LYS	A	244	11.232	68.728	25.122	1.00	11.69	6
1933	CG	LYS	A	244	10.924	70.142	24.580	1.00	14.16	6
1934	CD	LYS	A	244	9.398	70.400	24.483	1.00	13.14	6
1935	CE	LYS	A	244	9.129	71.869	24.133	1.00	13.82	6
1936	NZ	LYS	A	244	9.582	72.780	25.269	1.00	12.11	7
1937	N	LEU	A	245	12.213	67.068	27.593	1.00	11.20	7
1938	CA	LEU	A	245	12.730	65.762	28.039	1.00	11.98	6
1939	C	LEU	A	245	11.936	65.200	29.205	1.00	12.96	6
1940	O	LEU	A	245	11.665	63.996	29.219	1.00	12.16	8
1941	CB	LEU	A	245	14.221	65.961	28.461	1.00	10.88	6
1942	CG	LEU	A	245	15.091	66.282	27.206	1.00	15.60	6
1943	CD1	LEU	A	245	16.493	66.701	27.692	1.00	14.87	6
1944	CD2	LEU	A	245	15.227	65.052	26.282	1.00	19.24	6
1945	N	TYR	A	246	11.480	66.037	30.136	1.00	11.31	7
1946	CA	TYR	A	246	10.676	65.529	31.258	1.00	11.13	6
1947	C	TYR	A	246	9.294	65.090	30.770	1.00	11.59	6
1948	O	TYR	A	246	8.674	64.306	31.500	1.00	13.03	8
1949	CB	TYR	A	246	10.582	66.586	32.359	1.00	12.53	6
1950	CG	TYR	A	246	11.928	66.907	32.995	1.00	9.85	6
1951	CD1	TYR	A	246	12.882	65.941	33.210	1.00	11.79	6
1952	CD2	TYR	A	246	12.163	68.228	33.408	1.00	10.52	6
1953	CE1	TYR	A	246	14.103	66.245	33.830	1.00	11.71	6
1954	CE2	TYR	A	246	13.379	68.569	34.021	1.00	10.59	6
1955	CZ	TYR	A	246	14.319	67.562	34.208	1.00	11.60	6
1956	OH	TYR	A	246	15.536	67.856	34.816	1.00	11.49	8
1957	N	GLN	A	247	8.769	65.623	29.672	1.00	12.54	7
1958	CA	GLN	A	247	7.501	65.088	29.112	1.00	13.07	6
1959	C	GLN	A	247	7.677	63.690	28.587	1.00	13.75	6
1960	O	GLN	A	247	6.712	62.875	28.651	1.00	16.16	8
1961	CB	GLN	A	247	7.016	66.001	27.940	1.00	12.33	6
1962	CG	GLN	A	247	6.530	67.357	28.518	1.00	13.62	6
1963	CD	GLN	A	247	6.016	68.220	27.397	1.00	16.89	6
1964	OE1	GLN	A	247	5.355	67.699	26.462	1.00	18.89	8
1965	NE2	GLN	A	247	6.372	69.518	27.387	1.00	14.76	7
1966	N	LYS	A	248	8.881	63.349	28.162	1.00	14.17	7
1967	CA	LYS	A	248	9.163	61.979	27.702	1.00	15.18	6
1968	C	LYS	A	248	9.328	61.000	28.836	1.00	15.52	6
1969	O	LYS	A	248	8.839	59.868	28.746	1.00	16.45	8
1970	CB	ALYS	A	248	10.397	61.994	26.793	0.50	13.40	6
1971	CG	ALYS	A	248	10.116	62.793	25.528	0.50	14.14	6
1972	CD	ALYS	A	248	8.958	62.165	24.749	0.50	17.69	6
1973	CE	ALYS	A	248	8.449	63.068	23.657	0.50	18.41	6
1974	NZ	ALYS	A	248	7.682	62.378	22.577	0.50	25.68	6
1970	CB	BLYS	A	248	10.414	62.030	26.816	0.50	17.60	6

#	Atom	Res	Ch	Res#	X	Y	Z	Occ	B	El
1971	CG	BLYS	A	248	10.840	60.676	26.292	0.50	22.20	6
1972	CD	BLYS	A	248	11.561	60.755	24.977	0.50	29.85	6
1973	CE	BLYS	A	248	11.495	59.497	24.150	0.50	18.09	6
1974	NZ	BLYS	A	248	10.779	58.367	24.885	0.50	21.86	7
1975	N	LYS	A	249	10.131	61.400	29.830	1.00	12.97	7
1976	CA	LYS	A	249	10.424	60.442	30.917	1.00	14.04	6
1977	C	LYS	A	249	10.983	61.213	32.084	1.00	11.94	6
1978	O	LYS	A	249	11.520	62.316	31.876	1.00	13.26	8
1979	CB	LYS	A	249	11.514	59.423	30.462	1.00	15.09	6
1980	CG	LYS	A	249	11.674	58.358	31.542	1.00	15.17	6
1981	CD	LYS	A	249	12.552	57.175	31.147	1.00	23.08	6
1982	CE	LYS	A	249	12.451	56.111	32.249	1.00	27.89	6
1983	NZ	LYS	A	249	13.149	54.836	31.875	1.00	36.10	7
1984	N	ASP	A	250	10.870	60.678	33.288	1.00	11.41	7
1985	CA	ASP	A	250	11.406	61.336	34.494	1.00	12.04	6
1986	C	ASP	A	250	12.918	61.011	34.626	1.00	11.51	6
1987	O	ASP	A	250	13.348	60.354	35.546	1.00	13.89	8
1988	CB	ASP	A	250	10.638	60.891	35.728	1.00	11.54	6
1989	CG	ASP	A	250	10.618	59.405	36.030	1.00	15.08	6
1990	OD1	ASP	A	250	10.695	58.618	35.085	1.00	15.58	8
1991	OD2	ASP	A	250	10.448	59.027	37.224	1.00	15.73	8
1992	N	ILE	A	251	13.644	61.541	33.607	1.00	11.19	7
1993	CA	ILE	A	251	15.093	61.346	33.593	1.00	12.04	6
1994	C	ILE	A	251	15.777	62.413	34.467	1.00	12.14	6
1995	O	ILE	A	251	15.148	63.303	34.990	1.00	13.37	8
1996	CB	ILE	A	251	15.610	61.331	32.164	1.00	15.29	6
1997	CG1	ILE	A	251	14.988	62.477	31.361	1.00	23.77	6
1998	CG2	ILE	A	251	15.204	60.064	31.402	1.00	16.73	6
1999	CD1	ILE	A	251	15.645	63.784	31.503	1.00	25.03	6
2000	N	PHE	A	252	17.122	62.248	34.639	1.00	9.99	7
2001	CA	PHE	A	252	17.885	63.182	35.420	1.00	9.59	6
2002	C	PHE	A	252	18.550	64.105	34.501	1.00	11.26	6
2003	O	PHE	A	252	19.347	63.587	33.596	1.00	12.38	8
2004	CB	PHE	A	252	18.916	62.372	36.275	1.00	13.00	6
2005	CG	PHE	A	252	19.748	63.290	37.145	1.00	11.27	6
2006	CD1	PHE	A	252	19.182	63.951	38.241	1.00	12.24	6
2007	CD2	PHE	A	252	21.094	63.494	36.838	1.00	12.52	6
2008	CE1	PHE	A	252	19.907	64.802	39.029	1.00	11.05	6
2009	CE2	PHE	A	252	21.843	64.237	37.664	1.00	11.26	6
2010	CZ	PHE	A	252	21.263	65.031	38.750	1.00	10.84	6
2011	N	LEU	A	253	18.740	65.389	34.772	1.00	12.09	7
2012	CA	LEU	A	253	19.592	66.278	33.936	1.00	9.30	6
2013	C	LEU	A	253	20.498	67.083	34.913	1.00	9.94	6
2014	O	LEU	A	253	20.094	67.647	35.914	1.00	11.26	8
2015	CB	LEU	A	253	18.694	67.378	33.282	1.00	8.72	6
2016	CG	LEU	A	253	17.749	66.766	32.216	1.00	10.91	6
2017	CD1	LEU	A	253	16.881	67.891	31.562	1.00	10.35	6
2018	CD2	LEU	A	253	18.508	66.051	31.089	1.00	12.13	6
2019	N	VAL	A	254	21.761	67.179	34.422	1.00	11.23	7
2020	CA	VAL	A	254	22.750	68.015	35.077	1.00	11.09	6
2021	C	VAL	A	254	23.423	68.807	33.921	1.00	10.26	6
2022	O	VAL	A	254	23.707	68.196	32.892	1.00	10.32	8
2023	CB	VAL	A	254	23.722	67.184	32.765	1.00	9.61	6
2024	CG1	VAL	A	254	24.552	66.161	35.947	1.00	10.99	6
2025	CG2	VAL	A	254	24.688	68.131	35.161	1.00	11.22	6
2026	N	GLY	A	255	23.762	70.047	36.685	1.00	10.96	7
2027	CA	GLY	A	255	24.492	70.879	34.194	1.00	11.43	6
2028	C	GLY	A	255	25.877	71.193	33.243	1.00	11.23	6
2029	O	GLY	A	255	26.095	71.404	33.731	1.00	10.31	8
2030	N	GLU	A	256	26.828	71.308	34.967	1.00	10.72	7
2031	CA	GLU	A	256	28.159	71.786	33.187	1.00	10.84	6
2032	C	GLU	A	256	28.236	73.315	33.013	1.00	11.33	6
2033	O	GLU	A	256	28.295	73.820	31.871	1.00	10.36	8
2034	CB	GLU	A	256	29.172	71.178	32.167	1.00	10.52	6
2035	CG	GLU	A	256	30.603	71.617	32.605	1.00	10.70	6
2036	CD	GLU	A	256	31.442	72.059	31.414	1.00	11.32	6
2037	OE1	GLU	A	256	30.925	72.467	30.347	1.00	10.78	8
2038	OE2	GLU	A	256	32.696	71.998	31.517	1.00	11.88	8
2039	N	TRP	A	257	28.173	74.038	34.120	1.00	9.79	7
2040	CA	TRP	A	257	28.409	74.158	34.158	1.00	9.57	6
2041	C	TRP	A	257	29.798	75.480	34.799	1.00	10.84	6
2042	O	TRP	A	257	29.908	75.608	36.037	1.00	10.82	8
2043	CB	TRP	A	257	27.301	75.560	34.997	1.00	9.83	6
2044	CG	TRP	A	257	27.449	76.175	34.854	1.00	8.59	6
2045	CD1	TRP	A	257	28.566	77.700	35.196	1.00	12.90	6
2046	CD2	TRP	A	257	26.431	78.442	34.412	1.00	10.35	6
2047	NE1	TRP	A	257	28.295	78.592	34.937	1.00	12.55	7
2048	CE2	TRP	A	257	27.020	79.753	34.459	1.00	12.29	6
2049	CE3	TRP	A	257	25.124	79.891	33.946	1.00	13.94	6
2050	CZ2	TRP	A	257	26.317	78.425	34.055	1.00	10.73	6
2051	CZ3	TRP	A	257	24.385	81.041	33.527	1.00	13.32	6
2052	CH2	TRP	A	257	25.026	79.544	33.599	1.00	11.64	6
2053	N	TYR	A	258	30.831	80.793	33.986	1.00	11.79	7
2054	CA	TYR	A	258	32.211	75.725	34.524	1.00	11.80	6

2055	C	TYR	A	258	32.527	76.859	35.355	1.00	12.06	6
2056	O	TYR	A	258	32.370	77.988	34.869	1.00	13.57	8
2057	CB	TYR	A	258	33.189	75.508	33.335	1.00	11.31	6
2058	CG	TYR	A	258	34.562	74.978	33.723	1.00	11.97	6
2059	CD1	TYR	A	258	35.452	75.687	34.524	1.00	13.71	6
2060	CD2	TYR	A	258	34.932	73.734	33.234	1.00	11.78	6
2061	CE1	TYR	A	258	36.707	75.160	34.859	1.00	15.77	6
2062	CE2	TYR	A	258	36.179	73.226	33.564	1.00	13.38	6
2063	CZ	TYR	A	258	37.037	73.908	34.376	1.00	16.44	6
2064	OH	TYR	A	258	38.277	73.304	34.631	1.00	20.59	8
2065	N	GLY	A	259	33.030	76.605	36.550	1.00	9.86	7
2066	CA	GLY	A	259	33.584	77.705	37.361	1.00	12.24	6
2067	C	GLY	A	259	32.510	78.690	37.891	1.00	11.55	6
2068	O	GLY	A	259	31.417	78.276	38.227	1.00	12.61	8
2069	N	ASP	A	260	32.921	79.963	37.891	1.00	11.45	7
2070	CA	ASP	A	260	32.064	81.004	38.486	1.00	10.59	6
2071	C	ASP	A	260	31.718	80.614	39.947	1.00	12.08	6
2072	O	ASP	A	260	30.554	80.528	40.317	1.00	11.34	8
2073	CB	ASP	A	260	30.792	81.286	37.665	1.00	13.01	6
2074	CG	ASP	A	260	31.160	81.807	36.269	1.00	16.43	6
2075	OD1	ASP	A	260	32.136	82.593	36.148	1.00	13.58	8
2076	OD2	ASP	A	260	30.486	81.502	35.274	1.00	14.47	8
2077	N	ASP	A	261	32.812	80.528	40.708	1.00	10.89	7
2078	CA	ASP	A	261	32.709	80.239	42.158	1.00	13.96	6
2079	C	ASP	A	261	32.059	81.369	42.936	1.00	12.00	6
2080	O	ASP	A	261	31.920	82.528	42.502	1.00	13.08	8
2081	CB	ASP	A	261	34.125	79.875	42.610	1.00	17.34	6
2082	CG	ASP	A	261	34.615	78.518	42.074	1.00	18.47	6
2083	OD1	ASP	A	261	33.990	77.881	41.181	1.00	22.23	8
2084	OD2	ASP	A	261	35.642	78.035	42.569	1.00	20.46	8
2085	N	PRO	A	262	31.751	81.124	44.230	1.00	11.36	7
2086	CA	PRO	A	262	31.155	82.197	45.033	1.00	13.50	6
2087	C	PRO	A	262	32.085	83.428	45.042	1.00	15.27	6
2088	O	PRO	A	262	33.325	83.254	45.078	1.00	17.80	8
2089	CB	PRO	A	262	30.973	81.583	46.445	1.00	14.74	6
2090	CG	PRO	A	262	30.818	80.086	46.071	1.00	13.76	6
2091	CD	PRO	A	262	31.873	79.856	44.938	1.00	11.66	6
2092	N	GLY	A	263	31.385	84.560	45.069	1.00	20.11	7
2093	CA	GLY	A	263	32.178	85.810	45.091	1.00	20.82	6
2094	C	GLY	A	263	32.444	86.343	43.693	1.00	22.93	6
2095	O	GLY	A	263	33.030	87.433	43.565	1.00	26.12	8
2096	N	THR	A	264	32.204	85.581	42.640	1.00	16.30	7
2097	CA	THR	A	264	32.579	85.990	41.270	1.00	15.29	6
2098	C	THR	A	264	31.378	86.647	40.599	1.00	14.14	6
2099	O	THR	A	264	30.225	86.490	41.019	1.00	15.14	8
2100	CB	THR	A	264	33.154	84.890	40.377	1.00	17.00	6
2101	OG1	THR	A	264	32.185	83.864	40.214	1.00	17.20	8
2102	CG2	THR	A	264	34.455	84.302	40.998	1.00	18.53	6
2103	N	ALA	A	265	31.684	87.407	39.536	1.00	14.47	7
2104	CA	ALA	A	265	30.579	88.206	38.957	1.00	16.68	6
2105	C	ALA	A	265	29.455	87.387	38.348	1.00	14.50	6
2106	O	ALA	A	265	28.315	87.921	38.316	1.00	15.71	8
2107	CB	ALA	A	265	31.153	89.034	37.793	1.00	18.93	6
2108	N	ASN	A	266	29.808	86.186	37.837	1.00	11.26	7
2109	CA	ASN	A	266	28.739	85.402	37.199	1.00	12.93	6
2110	C	ASN	A	266	28.140	84.338	38.110	1.00	13.52	6
2111	O	ASN	A	266	27.364	83.484	37.632	1.00	11.82	8
2112	CB	ASN	A	266	29.289	84.792	35.855	1.00	15.03	6
2113	CG	ASN	A	266	29.632	85.944	34.889	1.00	15.26	6
2114	OD1	ASN	A	266	28.938	86.955	34.822	1.00	13.24	8
2115	ND2	ASN	A	266	30.698	85.797	34.146	1.00	14.64	7
2116	N	HIS	A	267	28.621	84.306	39.353	1.00	12.23	7
2117	CA	HIS	A	267	28.106	83.276	40.271	1.00	13.10	6
2118	C	HIS	A	267	26.596	83.276	40.469	1.00	13.44	6
2119	O	HIS	A	267	25.999	82.182	40.399	1.00	12.11	8
2120	CB	HIS	A	267	28.852	83.439	41.616	1.00	10.82	6
2121	CG	HIS	A	267	28.469	82.306	42.563	1.00	12.37	6
2122	ND1	HIS	A	267	28.877	81.020	42.410	1.00	11.65	7
2123	CD2	HIS	A	267	27.637	82.360	43.640	1.00	15.29	6
2124	CE1	HIS	A	267	28.355	80.278	43.375	1.00	12.97	6
2125	NE2	HIS	A	267	27.608	81.080	44.153	1.00	11.63	7
2126	N	LEU	A	268	26.001	84.455	40.988	1.00	10.81	7
2127	CA	LEU	A	268	24.548	84.430	40.726	1.00	13.40	6
2128	C	LEU	A	268	23.766	83.997	39.768	1.00	11.02	6
2129	O	LEU	A	268	22.745	83.268	39.957	1.00	12.74	8
2130	CB	LEU	A	268	24.169	85.889	41.374	1.00	14.95	6
2131	CG	LEU	A	268	22.599	86.052	41.471	1.00	19.99	6
2132	CD1	LEU	A	268	22.040	85.151	42.563	1.00	24.25	6
2133	CD2	LEU	A	268	22.298	87.536	41.677	1.00	25.18	6
2134	N	GLU	A	269	24.210	84.274	38.549	1.00	11.51	7
2135	CA	GLU	A	269	23.495	83.760	37.381	1.00	11.81	6
2136	C	GLU	A	269	23.525	82.221	37.386	1.00	10.11	6
2137	O	GLU	A	269	22.512	81.596	37.067	1.00	11.07	8
2138	CB	GLU	A	269	24.190	84.360	36.122	1.00	13.26	6

2139	CG	GLU	A	269	23.490	83.803	34.867	1.00	11.45	6
2140	CD	GLU	A	269	24.122	84.417	33.582	1.00	13.11	6
2141	OE1	GLU	A	269	24.906	85.369	33.643	1.00	12.92	8
2142	OE2	GLU	A	269	23.720	83.906	32.525	1.00	12.17	8
2143	N	LYS	A	270	24.678	81.608	37.720	1.00	9.00	7
2144	CA	LYS	A	270	24.754	80.156	37.787	1.00	9.72	6
2145	C	LYS	A	270	23.806	79.578	38.862	1.00	10.98	6
2146	O	LYS	A	270	23.079	78.613	38.609	1.00	10.64	8
2147	CB	LYS	A	270	26.221	79.709	38.063	1.00	10.49	6
2148	CG	LYS	A	270	26.259	78.174	38.299	1.00	10.05	6
2149	CD	LYS	A	270	27.725	77.682	38.502	1.00	8.77	6
2150	CE	LYS	A	270	28.236	78.181	39.869	1.00	9.57	6
2151	NZ	LYS	A	270	29.600	77.515	40.161	1.00	11.56	7
2152	N	VAL	A	271	23.765	80.218	40.038	1.00	11.90	7
2153	CA	VAL	A	271	22.847	79.712	41.105	1.00	10.97	6
2154	C	VAL	A	271	21.392	79.835	40.608	1.00	10.81	6
2155	O	VAL	A	271	20.583	78.951	40.827	1.00	11.67	8
2156	CB	VAL	A	271	23.044	80.608	42.346	1.00	13.53	6
2157	CG1	VAL	A	271	22.016	80.248	43.435	1.00	12.11	6
2158	CG2	VAL	A	271	24.455	80.307	42.888	1.00	12.46	6
2159	N	ARG	A	272	21.064	80.978	39.983	1.00	11.20	7
2160	CA	ARG	A	272	19.668	81.127	39.458	1.00	11.53	6
2161	C	ARG	A	272	19.328	80.041	38.439	1.00	10.29	6
2162	O	ARG	A	272	18.208	79.506	38.389	1.00	11.34	8
2163	CB	ARG	A	272	19.462	82.540	38.859	1.00	9.74	6
2164	CG	ARG	A	272	19.220	83.552	40.027	1.00	11.97	6
2165	CD	ARG	A	272	19.405	85.000	39.483	1.00	10.96	6
2166	NE	ARG	A	272	18.600	85.394	38.306	1.00	12.10	7
2167	CZ	ARG	A	272	17.297	85.759	38.377	1.00	14.61	6
2168	NH1	ARG	A	272	16.541	85.765	39.493	1.00	12.14	7
2169	NH2	ARG	A	272	16.734	86.132	37.232	1.00	12.07	7
2170	N	TYR	A	273	20.329	79.790	37.553	1.00	10.33	7
2171	CA	TYR	A	273	20.116	78.669	36.589	1.00	11.57	6
2172	C	TYR	A	273	19.882	77.355	37.328	1.00	9.14	6
2173	O	TYR	A	273	18.931	76.590	36.993	1.00	12.29	8
2174	CB	TYR	A	273	21.415	78.569	35.742	1.00	9.07	6
2175	CG	TYR	A	273	21.388	77.445	34.693	1.00	8.78	6
2176	CD1	TYR	A	273	21.708	76.144	35.110	1.00	11.52	6
2177	CD2	TYR	A	273	21.054	77.630	33.363	1.00	9.55	6
2178	CE1	TYR	A	273	21.663	75.071	34.218	1.00	9.62	6
2179	CE2	TYR	A	273	21.087	76.577	32.482	1.00	8.45	6
2180	CZ	TYR	A	273	21.366	75.301	32.885	1.00	10.47	6
2181	OH	TYR	A	273	21.365	74.258	31.956	1.00	11.35	8
2182	N	ALA	A	274	20.678	77.001	38.308	1.00	10.68	7
2183	CA	ALA	A	274	20.449	75.697	39.008	1.00	10.66	6
2184	C	ALA	A	274	19.062	75.725	39.677	1.00	10.79	6
2185	O	ALA	A	274	18.478	74.649	39.796	1.00	11.53	8
2186	CB	ALA	A	274	21.539	75.560	40.092	1.00	12.82	7
2187	N	ASN	A	275	18.692	76.887	40.217	1.00	11.15	6
2188	CA	ASN	A	275	17.443	76.882	40.998	1.00	11.26	6
2189	C	ASN	A	275	16.199	77.051	40.131	1.00	10.88	8
2190	O	ASN	A	275	15.082	76.705	40.606	1.00	11.83	6
2191	CB	ASN	A	275	17.490	78.056	42.004	1.00	11.31	6
2192	CG	ASN	A	275	18.495	77.768	43.107	1.00	11.97	8
2193	OD1	ASN	A	275	18.987	76.661	43.214	1.00	12.55	7
2194	ND2	ASN	A	275	18.759	78.816	43.934	1.00	12.55	7
2195	N	ASN	A	276	16.322	77.453	38.842	1.00	11.20	6
2196	CA	ASN	A	276	15.129	77.705	38.045	1.00	12.25	6
2197	C	ASN	A	276	15.023	76.992	36.720	1.00	14.18	8
2198	O	ASN	A	276	13.932	76.820	36.165	1.00	11.89	6
2199	CB	ASN	A	276	15.134	79.197	37.632	1.00	9.94	6
2200	CG	ASN	A	276	14.629	80.023	38.806	1.00	14.75	6
2201	OD1	ASN	A	276	13.365	80.089	38.826	1.00	16.83	8
2202	ND2	ASN	A	276	15.551	80.529	39.612	1.00	16.12	7
2203	N	SER	A	277	16.151	76.504	36.173	1.00	10.52	6
2204	CA	SER	A	277	16.064	75.974	34.811	1.00	11.82	6
2205	C	SER	A	277	15.497	74.585	34.678	1.00	11.16	6
2206	O	SER	A	277	15.204	74.178	33.545	1.00	12.16	8
2207	CB	SER	A	277	17.502	75.911	34.204	1.00	11.66	6
2208	OG	SER	A	277	18.257	74.877	34.872	1.00	12.14	8
2209	N	GLY	A	278	15.371	73.851	35.795	1.00	11.36	7
2210	CA	GLY	A	278	15.051	72.411	35.697	1.00	11.40	6
2211	C	GLY	A	278	16.263	71.565	35.375	1.00	11.97	6
2212	O	GLY	A	278	16.115	70.348	35.279	1.00	13.61	8
2213	N	VAL	A	279	17.442	72.180	35.342	1.00	12.14	7
2214	CA	VAL	A	279	18.673	71.427	35.090	1.00	10.48	6
2215	C	VAL	A	279	19.552	71.645	36.348	1.00	11.31	6
2216	O	VAL	A	279	19.814	72.782	36.732	1.00	12.60	8
2217	CB	VAL	A	279	19.362	71.989	33.830	1.00	11.56	6
2218	CG1	VAL	A	279	20.628	71.191	33.543	1.00	10.87	6
2219	CG2	VAL	A	279	18.429	71.824	32.592	1.00	13.38	6
2220	N	ASN	A	280	20.053	70.556	36.919	1.00	10.03	7
2221	CA	ASN	A	280	20.918	70.714	38.122	1.00	10.49	6
2222	C	ASN	A	280	22.294	71.095	37.572	1.00	13.05	6

70

Atom#	Atom	Res	Chain	ResNum	X	Y	Z	Occ	B	Elem
2223	O	ASN	A	280	22.506	71.116	36.365	1.00	12.57	8
2224	CB	ASN	A	280	20.968	69.368	38.864	1.00	11.26	6
2225	CG	ASN	A	280	19.492	69.056	39.263	1.00	9.85	6
2226	OD1	ASN	A	280	18.839	69.805	39.990	1.00	10.91	8
2227	ND2	ASN	A	280	18.974	67.913	38.787	1.00	10.65	7
2228	N	VAL	A	281	23.262	71.312	38.492	1.00	11.45	7
2229	CA	VAL	A	281	24.622	71.621	37.977	1.00	9.71	6
2230	C	VAL	A	281	25.671	70.782	38.660	1.00	11.24	6
2231	O	VAL	A	281	25.581	70.318	39.799	1.00	11.15	8
2232	CB	VAL	A	281	25.013	73.104	38.187	1.00	9.82	6
2233	CG1	VAL	A	281	24.019	74.024	37.397	1.00	10.47	6
2234	CG2	VAL	A	281	25.038	73.569	39.638	1.00	11.52	6
2235	N	LEU	A	282	26.786	70.675	37.904	1.00	9.31	7
2236	CA	LEU	A	282	28.031	70.087	38.460	1.00	10.27	6
2237	C	LEU	A	282	28.631	71.063	39.476	1.00	11.17	6
2238	O	LEU	A	282	28.577	72.295	39.298	1.00	12.65	8
2239	CB	LEU	A	282	29.022	69.883	37.283	1.00	10.13	6
2240	CG	LEU	A	282	28.650	68.558	36.538	1.00	11.97	6
2241	CD1	LEU	A	282	29.159	68.726	35.102	1.00	16.51	6
2242	CD2	LEU	A	282	29.314	67.351	37.196	1.00	12.93	6
2243	N	ASP	A	283	29.128	70.458	40.581	1.00	9.33	7
2244	CA	ASP	A	283	29.503	71.304	41.764	1.00	11.76	6
2245	C	ASP	A	283	30.935	71.778	41.696	1.00	11.31	6
2246	O	ASP	A	283	31.877	71.275	42.316	1.00	11.22	8
2247	CB	ASP	A	283	29.212	70.451	43.014	1.00	8.73	6
2248	CG	ASP	A	283	29.282	71.343	44.271	1.00	11.60	6
2249	OD1	ASP	A	283	29.655	72.526	44.259	1.00	10.06	8
2250	OD2	ASP	A	283	28.866	70.758	45.290	1.00	11.95	8
2251	N	PHE	A	284	31.108	72.815	40.834	1.00	10.41	7
2252	CA	PHE	A	284	32.439	73.424	40.731	1.00	10.33	6
2253	C	PHE	A	284	32.746	74.271	41.966	1.00	11.21	6
2254	O	PHE	A	284	33.941	74.331	42.313	1.00	12.52	8
2255	CB	PHE	A	284	32.509	74.409	39.517	1.00	10.51	6
2256	CG	PHE	A	284	32.750	73.575	38.227	1.00	10.01	6
2257	CD1	PHE	A	284	31.725	72.991	37.520	1.00	11.07	6
2258	CD2	PHE	A	284	34.073	73.416	37.765	1.00	11.57	6
2259	CE1	PHE	A	284	31.925	72.232	36.386	1.00	12.29	6
2260	CE2	PHE	A	284	34.292	72.665	36.596	1.00	11.45	6
2261	CZ	PHE	A	284	33.234	72.085	35.903	1.00	9.72	6
2262	N	ASP	A	285	31.729	74.753	42.687	1.00	9.13	7
2263	CA	ASP	A	285	32.061	75.575	43.882	1.00	10.61	6
2264	C	ASP	A	285	32.799	74.670	44.885	1.00	10.95	6
2265	O	ASP	A	285	33.806	75.092	45.468	1.00	11.00	8
2266	CB	ASP	A	285	30.773	76.068	44.545	1.00	9.55	6
2267	CG	ASP	A	285	29.979	77.042	43.692	1.00	10.92	6
2268	OD1	ASP	A	285	30.499	77.631	42.710	1.00	12.04	8
2269	OD2	ASP	A	285	28.773	77.214	44.043	1.00	12.59	8
2270	N	LEU	A	286	32.248	73.464	45.128	1.00	10.30	7
2271	CA	LEU	A	286	32.959	72.569	46.061	1.00	9.08	6
2272	C	LEU	A	286	34.219	71.940	45.505	1.00	9.26	6
2273	O	LEU	A	286	35.179	71.757	46.238	1.00	11.04	8
2274	CB	LEU	A	286	31.968	71.419	46.481	1.00	9.12	6
2275	CG	LEU	A	286	32.367	70.697	47.790	1.00	10.68	6
2276	CD1	LEU	A	286	32.357	71.580	49.023	1.00	15.93	6
2277	CD2	LEU	A	286	31.384	69.510	47.999	1.00	13.68	6
2278	N	ASN	A	287	34.226	71.683	44.168	1.00	9.32	7
2279	CA	ASN	A	287	35.441	71.091	43.586	1.00	10.65	6
2280	C	ASN	A	287	36.687	71.937	43.872	1.00	9.75	6
2281	O	ASN	A	287	37.751	71.401	44.184	1.00	10.35	8
2282	CB	ASN	A	287	35.269	71.000	42.039	1.00	8.74	6
2283	CG	ASN	A	287	36.571	70.507	41.392	1.00	11.79	6
2284	OD1	ASN	A	287	37.171	71.318	40.654	1.00	11.65	8
2285	ND2	ASN	A	287	36.866	69.245	41.597	1.00	9.92	7
2286	N	THR	A	288	36.588	73.275	43.719	1.00	9.52	7
2287	CA	THR	A	288	37.769	74.106	43.907	1.00	11.58	6
2288	C	THR	A	288	38.367	73.906	45.292	1.00	10.74	6
2289	O	THR	A	288	39.593	73.741	45.486	1.00	13.25	8
2290	CB	THR	A	288	37.365	75.596	43.745	1.00	15.82	6
2291	OG1	THR	A	288	36.906	75.753	42.397	1.00	16.55	8
2292	CG2	THR	A	288	38.641	76.459	43.984	1.00	14.43	6
2293	N	VAL	A	289	37.440	73.847	46.268	1.00	8.64	7
2294	CA	VAL	A	289	37.915	73.732	47.655	1.00	8.82	6
2295	C	VAL	A	289	38.401	72.341	47.995	1.00	11.04	6
2296	O	VAL	A	289	39.438	72.184	48.642	1.00	11.67	8
2297	CB	VAL	A	289	36.780	74.237	48.605	1.00	11.02	6
2298	CG1	VAL	A	289	37.297	74.258	50.051	1.00	10.38	6
2299	CG2	VAL	A	289	36.323	75.632	48.191	1.00	11.61	6
2300	N	ILE	A	290	37.764	73.287	47.430	1.00	9.86	7
2301	CA	ILE	A	290	38.262	69.911	47.642	1.00	8.30	6
2302	C	ILE	A	290	39.715	69.830	47.144	1.00	12.22	6
2303	O	ILE	A	290	40.554	69.251	47.806	1.00	10.89	8
2304	CB	ILE	A	290	37.342	68.979	46.851	1.00	9.93	6
2305	CG1	ILE	A	290	36.004	68.833	47.591	1.00	11.46	6
2306	CG2	ILE	A	290	38.025	67.579	46.690	1.00	11.35	6

#	Atom	Res	Chain	Seq	X	Y	Z	Occ	B	n
2307	CD1	ILE	A	290	34.964	68.108	46.700	1.00	11.71	6
2308	N	ARG	A	291	39.969	70.386	45.940	1.00	11.21	7
2309	CA	ARG	A	291	41.352	70.246	45.455	1.00	10.67	6
2310	C	ARG	A	291	42.343	71.071	46.293	1.00	9.59	6
2311	O	ARG	A	291	43.481	70.627	46.380	1.00	12.39	8
2312	CB	ARG	A	291	41.308	70.770	43.991	1.00	10.96	6
2313	CG	ARG	A	291	40.602	69.717	43.110	1.00	11.39	6
2314	CD	ARG	A	291	40.480	70.173	41.647	1.00	11.11	6
2315	NE	ARG	A	291	41.729	70.343	40.907	1.00	11.15	7
2316	CZ	ARG	A	291	42.225	69.361	40.109	1.00	12.12	6
2317	NH1	ARG	A	291	41.706	68.126	40.068	1.00	11.32	7
2318	NH2	ARG	A	291	43.322	69.593	39.363	1.00	12.38	7
2319	N	ASN	A	292	41.972	72.204	46.801	1.00	12.65	7
2320	CA	ASN	A	292	42.894	73.001	47.653	1.00	11.32	6
2321	C	ASN	A	292	42.983	72.376	49.051	1.00	11.68	6
2322	O	ASN	A	292	44.074	72.565	49.634	1.00	11.50	8
2323	CB	ASN	A	292	42.408	74.471	47.750	1.00	11.01	6
2324	CG	ASN	A	292	42.593	75.238	46.379	1.00	14.57	6
2325	OD1	ASN	A	292	43.466	74.831	45.665	1.00	19.84	8
2326	ND2	ASN	A	292	41.735	76.242	46.267	1.00	18.55	7
2327	N	VAL	A	293	42.009	71.574	49.542	1.00	10.40	7
2328	CA	VAL	A	293	42.157	71.075	50.940	1.00	10.24	6
2329	C	VAL	A	293	42.861	69.743	50.969	1.00	11.21	6
2330	O	VAL	A	293	43.748	69.520	51.783	1.00	11.98	8
2331	CB	VAL	A	293	40.703	70.961	51.499	1.00	10.95	6
2332	CG1	VAL	A	293	40.707	70.175	52.820	1.00	12.97	6
2333	CG2	VAL	A	293	42.479	72.364	51.736	1.00	13.88	6
2334	N	PHE	A	294	43.106	68.858	50.033	1.00	11.12	7
2335	CA	PHE	A	294	44.255	67.537	49.960	1.00	11.38	6
2336	C	PHE	A	294	45.095	67.464	48.964	1.00	11.51	6
2337	O	PHE	A	294	42.063	66.540	49.093	1.00	11.86	8
2338	CB	PHE	A	294	40.936	66.455	49.553	1.00	11.70	6
2339	CG	PHE	A	294	41.178	66.334	50.584	1.00	11.57	6
2340	CD1	PHE	A	294	39.707	65.683	51.808	1.00	11.48	6
2341	CD2	PHE	A	294	40.133	66.865	50.295	1.00	13.60	7
2342	CE1	PHE	A	294	38.662	65.579	52.724	1.00	12.41	6
2343	CE2	PHE	A	294	38.880	66.762	51.232	1.00	13.36	6
2344	CZ	PHE	A	294	44.295	66.104	52.463	1.00	13.22	7
2345	N	GLY	A	295	45.328	68.355	47.976	1.00	11.76	6
2346	CA	GLY	A	295	46.504	68.234	46.907	1.00	12.36	6
2347	C	GLY	A	295	47.547	69.187	47.187	1.00	10.85	6
2348	O	GLY	A	295		68.676	47.612	1.00	13.29	8
2349	N	THR	A	296	46.271	70.499	46.990	1.00	10.00	7
2350	CA	THR	A	296	47.441	71.394	47.123	1.00	11.19	6
2351	C	THR	A	296	47.675	71.894	48.549	1.00	13.50	6
2352	O	THR	A	296	48.755	72.436	48.874	1.00	12.22	8
2353	CB	THR	A	296	47.162	72.647	46.253	1.00	11.65	6
2354	OG1	THR	A	296	46.005	73.276	46.812	1.00	15.42	8
2355	CG2	THR	A	296	46.976	72.308	44.770	1.00	13.16	6
2356	N	PHE	A	297	46.760	71.647	49.472	1.00	11.24	7
2357	CA	PHE	A	297	46.917	72.026	50.881	1.00	12.10	6
2358	C	PHE	A	297	47.104	73.543	50.984	1.00	17.18	6
2359	O	PHE	A	297	47.719	74.035	51.952	1.00	17.61	8
2360	CB	PHE	A	297	48.052	71.271	51.610	1.00	11.80	6
2361	CG	PHE	A	297	47.630	69.924	52.199	1.00	14.38	6
2362	CD1	PHE	A	297	47.153	68.917	51.385	1.00	13.09	6
2363	CD2	PHE	A	297	47.760	69.711	53.565	1.00	11.70	6
2364	CE1	PHE	A	297	46.821	67.656	51.854	1.00	12.95	6
2365	CE2	PHE	A	297	47.438	68.444	54.064	1.00	12.92	6
2366	CZ	PHE	A	297	46.948	67.450	53.231	1.00	11.87	7
2367	N	THR	A	298	46.411	74.322	50.183	1.00	13.51	6
2368	CA	THR	A	298	46.398	75.773	50.270	1.00	13.52	6
2369	C	THR	A	298	45.131	76.303	50.925	1.00	13.29	6
2370	O	THR	A	298	45.018	77.505	51.220	1.00	15.14	8
2371	CB	THR	A	298	46.488	76.437	48.870	1.00	13.68	6
2372	OG1	THR	A	298	45.446	75.982	48.035	1.00	12.95	8
2373	CG2	THR	A	298	47.869	76.059	48.250	1.00	15.45	6
2374	N	GLN	A	299	44.162	75.400	51.194	1.00	10.17	7
2375	CA	GLN	A	299	43.009	75.785	51.975	1.00	13.00	6
2376	C	GLN	A	299	42.852	74.738	53.114	1.00	14.72	6
2377	O	GLN	A	299	43.624	73.753	53.094	1.00	14.92	8
2378	CB	GLN	A	299	41.654	75.808	51.214	1.00	11.79	6
2379	CG	GLN	A	299	41.692	77.072	50.299	1.00	14.80	6
2380	CD	GLN	A	299	40.301	77.320	49.668	1.00	16.23	6
2381	OE1	GLN	A	299	39.959	76.525	48.827	1.00	15.11	8
2382	NE2	GLN	A	299	39.635	78.376	50.107	1.00	17.75	7
2383	N	THR	A	300	42.031	75.020	54.106	1.00	13.71	7
2384	CA	THR	A	300	41.924	74.106	55.252	1.00	11.61	6
2385	C	THR	A	300	40.518	73.485	55.355	1.00	14.08	6
2386	O	THR	A	300	39.580	73.874	54.644	1.00	11.91	8
2387	CB	THR	A	300	42.152	74.850	56.601	1.00	15.34	6
2388	OG1	THR	A	300	41.116	75.771	56.804	1.00	15.70	8
2389	CG2	THR	A	300	43.511	75.587	56.464	1.00	16.98	6
2390	N	MET	A	301	40.337	72.672	56.443	1.00	11.88	7

2391	CA	MET	A	301	38.977	72.123	56.650	1.00	12.22	6
2392	C	MET	A	301	37.972	73.216	56.994	1.00	12.44	6
2393	O	MET	A	301	36.791	73.075	56.691	1.00	11.03	8
2394	CB	MET	A	301	38.935	71.065	57.800	1.00	12.80	6
2395	CG	MET	A	301	39.707	69.787	57.393	1.00	11.05	6
2396	SD	MET	A	301	39.027	69.014	55.895	1.00	12.28	16
2397	CE	MET	A	301	39.724	67.366	56.047	1.00	14.34	6
2398	N	TYR	A	302	38.408	74.374	57.555	1.00	12.00	7
2399	CA	TYR	A	302	37.462	75.453	57.759	1.00	10.79	6
2400	C	TYR	A	302	36.898	75.974	56.430	1.00	11.89	6
2401	O	TYR	A	302	35.694	76.181	56.325	1.00	12.65	8
2402	CB	TYR	A	302	38.131	76.626	58.505	1.00	10.20	6
2403	CG	TYR	A	302	38.409	76.271	59.983	1.00	11.51	6
2404	CD1	TYR	A	302	37.375	76.197	60.914	1.00	16.19	6
2405	CD2	TYR	A	302	39.739	76.057	60.344	1.00	17.97	6
2406	CE1	TYR	A	302	37.736	75.884	62.236	1.00	18.06	6
2407	CE2	TYR	A	302	40.062	75.723	61.683	1.00	15.93	6
2408	CZ	TYR	A	302	39.029	75.670	62.567	1.00	19.00	6
2409	OH	TYR	A	302	39.409	75.355	63.905	1.00	21.46	8
2410	N	ASP	A	303	37.763	76.026	55.418	1.00	13.06	7
2411	CA	ASP	A	303	37.256	76.434	54.104	1.00	12.54	6
2412	C	ASP	A	303	36.285	75.409	53.492	1.00	11.88	6
2413	O	ASP	A	303	35.330	75.782	52.785	1.00	13.46	8
2414	CB	ASP	A	303	38.419	76.666	53.133	1.00	10.93	6
2415	CG	ASP	A	303	39.386	77.734	53.654	1.00	14.78	6
2416	OD1	ASP	A	303	38.879	78.877	53.849	1.00	13.77	8
2417	OD2	ASP	A	303	40.573	77.435	53.832	1.00	13.10	8
2418	N	LEU	A	304	36.602	74.152	53.720	1.00	11.36	7
2419	CA	LEU	A	304	35.713	73.086	53.156	1.00	11.06	6
2420	C	LEU	A	304	34.348	73.181	53.841	1.00	12.95	6
2421	O	LEU	A	304	33.315	73.111	53.155	1.00	11.60	8
2422	CB	LEU	A	304	36.393	71.742	53.367	1.00	10.48	6
2423	CG	LEU	A	304	35.687	70.566	52.633	1.00	11.09	6
2424	CD1	LEU	A	304	35.737	70.797	51.112	1.00	12.57	6
2425	CD2	LEU	A	304	36.397	69.248	52.971	1.00	12.54	6
2426	N	ASN	A	305	34.290	73.255	55.180	1.00	11.76	7
2427	CA	ASN	A	305	32.999	73.422	55.887	1.00	12.74	6
2428	C	ASN	A	305	32.308	74.720	55.479	1.00	11.54	6
2429	O	ASN	A	305	31.100	74.673	55.259	1.00	13.50	8
2430	CB	ASN	A	305	33.331	73.448	57.405	1.00	12.22	6
2431	CG	ASN	A	305	32.014	73.551	58.210	1.00	12.84	6
2432	OD1	ASN	A	305	31.174	72.669	58.097	1.00	13.98	8
2433	ND2	ASN	A	305	31.910	74.651	58.967	1.00	17.24	7
2434	N	ASN	A	306	33.061	75.772	55.272	1.00	11.86	7
2435	CA	ASN	A	306	32.418	77.036	54.848	1.00	12.68	6
2436	C	ASN	A	306	31.740	76.830	53.499	1.00	13.44	6
2437	O	ASN	A	306	30.672	77.409	53.225	1.00	13.40	8
2438	CB	ASN	A	306	33.438	78.176	54.774	1.00	12.13	6
2439	CG	ASN	A	306	33.863	78.686	56.143	1.00	20.79	6
2440	OD1	ASN	A	306	33.275	78.499	57.233	1.00	23.19	8
2441	ND2	ASN	A	306	34.959	79.477	56.155	1.00	24.17	7
2442	N	MET	A	307	32.405	76.105	52.589	1.00	11.55	7
2443	CA	MET	A	307	31.750	75.915	51.254	1.00	11.38	6
2444	C	MET	A	307	30.590	74.957	51.272	1.00	12.48	6
2445	O	MET	A	307	29.639	75.156	50.512	1.00	13.10	8
2446	CB	MET	A	307	32.849	75.431	50.252	1.00	12.11	6
2447	CG	MET	A	307	31.759	75.462	48.767	1.00	12.14	6
2448	SD	MET	A	307	31.759	77.075	48.246	1.00	12.77	16
2449	CE	MET	A	307	33.282	78.010	48.324	1.00	13.86	6
2450	N	VAL	A	308	30.592	73.951	52.152	1.00	10.38	7
2451	CA	VAL	A	308	29.383	73.136	52.340	1.00	10.10	6
2452	C	VAL	A	308	28.272	74.049	52.822	1.00	11.64	6
2453	O	VAL	A	308	27.153	73.974	52.321	1.00	13.14	8
2454	CB	VAL	A	308	29.712	72.052	53.406	1.00	13.16	6
2455	CG1	VAL	A	308	28.388	71.356	53.770	1.00	14.05	6
2456	CG2	VAL	A	308	30.641	71.038	52.727	1.00	13.08	6
2457	N	ASN	A	309	28.541	74.952	53.784	1.00	12.04	7
2458	CA	ASN	A	309	27.479	75.848	54.264	1.00	12.56	6
2459	C	ASN	A	309	27.073	76.824	53.176	1.00	13.56	6
2460	O	ASN	A	309	25.875	77.017	53.030	1.00	15.38	8
2461	CB	ASN	A	309	27.997	76.680	55.470	1.00	14.00	6
2462	CG	ASN	A	309	28.109	75.761	56.684	1.00	20.56	6
2463	OD1	ASN	A	309	27.432	74.725	56.748	1.00	26.71	8
2464	ND2	ASN	A	309	28.967	76.208	57.595	1.00	21.18	7
2465	N	GLN	A	310	27.970	77.423	52.426	1.00	14.28	7
2466	CA	GLN	A	310	27.549	78.422	51.422	1.00	13.95	6
2467	C	GLN	A	310	26.734	77.754	50.319	1.00	13.50	6
2468	O	GLN	A	310	25.672	78.286	49.942	1.00	14.03	8
2469	CB	GLN	A	310	28.821	79.123	50.869	1.00	14.99	6
2470	CG	GLN	A	310	28.507	80.214	49.859	1.00	21.08	6
2471	CD	GLN	A	310	29.734	81.169	49.819	1.00	21.08	6
2472	OE1	GLN	A	310	30.875	80.762	50.019	1.00	25.61	8
2473	NE2	GLN	A	310	29.417	82.401	49.600	1.00	25.23	7
2474	N	THR	A	311	27.238	76.630	49.792	1.00	11.65	7

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2475  CA  THR A 311   26.432  75.983  48.711  1.00 11.52  6
2476  C   THR A 311   25.089  75.459  49.257  1.00 12.57  6
2477  O   THR A 311   24.039  75.572  48.599  1.00 13.81  8
2478  CB  THR A 311   27.143  74.754  48.142  1.00 12.52  6
2479  OG1 THR A 311   27.593  73.909  49.194  1.00 13.67  8
2480  CG2 THR A 311   28.426  75.250  47.375  1.00 14.32  6
2481  N   GLY A 312   25.096  75.107  50.561  1.00 11.69  7
2482  CA  GLY A 312   23.812  74.642  51.137  1.00 14.78  6
2483  C   GLY A 312   22.800  75.792  51.223  1.00 14.32  6
2484  O   GLY A 312   21.573  75.473  51.276  1.00 16.33  8
2485  N   ASN A 313   23.320  76.998  51.476  1.00 13.08  7
2486  CA  ASN A 313   22.392  78.108  51.535  1.00 13.20  6
2487  C   ASN A 313   21.980  78.575  50.150  1.00 14.89  6
2488  O   ASN A 313   20.827  79.097  50.014  1.00 20.75  8
2489  CB  ASN A 313   23.155  79.260  52.204  1.00 19.18  6
2490  CG  ASN A 313   23.210  79.024  53.718  1.00 30.91  6
2491  OD1 ASN A 313   22.384  78.310  54.281  1.00 30.04  8
2492  ND2 ASN A 313   24.152  79.691  54.360  1.00 31.03  7
2493  N   GLU A 314   22.802  78.394  49.111  1.00 12.83  7
2494  CA  GLU A 314   22.396  79.060  47.822  1.00 11.95  6
2495  C   GLU A 314   21.621  78.137  46.889  1.00 13.20  6
2496  O   GLU A 314   20.733  78.689  47.117  1.00 13.02  8
2497  CB  GLU A 314   23.753  79.394  47.842  1.00 10.17  6
2498  CG  GLU A 314   24.503  80.510  46.985  1.00 12.02  6
2499  CD  GLU A 314   25.715  80.987  46.232  1.00 15.94  6
2500  OE1 GLU A 314   26.285  80.206  47.284  1.00 15.22  8
2501  OE2 GLU A 314   26.164  82.100  46.892  1.00 24.00  8
2502  N   TYR A 315   21.992  76.840  45.885  1.00 11.76  7
2503  CA  TYR A 315   21.297  75.979  45.494  1.00 10.85  6
2504  C   TYR A 315   20.032  75.368  44.699  1.00 12.73  6
2505  O   TYR A 315   20.140  74.648  43.316  1.00 12.47  8
2506  CB  TYR A 315   22.265  74.838  45.297  1.00 12.39  6
2507  CG  TYR A 315   23.437  75.394  42.586  1.00 10.41  6
2508  CD1 TYR A 315   23.270  75.610  44.563  1.00 11.23  6
2509  CD2 TYR A 315   24.613  75.785  43.187  1.00 12.52  6
2510  CE1 TYR A 315   24.333  76.166  42.517  1.00 12.92  6
2511  CE2 TYR A 315   25.686  76.350  45.861  1.00 10.90  6
2512  CZ  TYR A 315   25.510  76.507  46.395  1.00 12.76  6
2513  OH  TYR A 315   26.595  77.056  46.454  1.00 13.74  8
2514  N   LYS A 316   18.895  75.634  45.861  1.00 11.54  7
2515  CA  LYS A 316   17.638  75.080  46.395  1.00 14.39  6
2516  C   LYS A 316   17.578  73.555  46.454  1.00 15.27  6

2517  O   LYS A 316   17.175  72.963  47.475  1.00 14.60  8
2518  CB  LYS A 316   16.496  75.595  45.485  1.00 14.77  6
2519  CG  LYS A 316   15.139  75.068  46.091  1.00 18.12  6
2520  CD  LYS A 316   13.982  75.856  45.565  1.00 22.24  6
2521  CE  LYS A 316   12.683  75.166  46.061  1.00 19.23  6
2522  NZ  LYS A 316   12.432  75.635  47.468  1.00 22.20  7
2523  N   TYR A 317   18.227  72.953  45.457  1.00 13.10  7
2524  CA  TYR A 317   18.316  71.483  45.445  1.00 11.39  6
2525  C   TYR A 317   19.805  71.123  45.506  1.00 10.31  6
2526  O   TYR A 317   20.410  70.491  44.637  1.00 12.52  8
2527  CB  TYR A 317   17.652  70.893  44.157  1.00 12.53  6
2528  CG  TYR A 317   16.221  71.387  44.004  1.00 13.61  6
2529  CD1 TYR A 317   15.272  70.932  44.915  1.00 14.65  6
2530  CD2 TYR A 317   15.800  72.280  43.030  1.00 14.34  6
2531  CE1 TYR A 317   13.938  71.366  44.811  1.00 13.78  6
2532  CE2 TYR A 317   14.511  72.767  42.890  1.00 12.10  6
2533  CZ  TYR A 317   13.604  72.246  43.832  1.00 15.19  6
2534  OH  TYR A 317   12.275  72.682  43.733  1.00 14.53  8
2535  N   LYS A 318   20.375  71.463  46.702  1.00 11.46  7
2536  CA  LYS A 318   21.821  71.130  46.865  1.00 10.64  6
2537  C   LYS A 318   22.053  69.651  46.865  1.00 10.21  6
2538  O   LYS A 318   23.133  69.145  46.529  1.00 10.36  8
2539  CB  LYS A 318   22.408  71.772  48.141  1.00  9.85  6
2540  CG  LYS A 318   21.843  71.211  49.451  1.00 13.57  6
2541  CD  LYS A 318   20.619  72.105  49.820  1.00 19.29  6
2542  CE  LYS A 318   20.309  72.064  51.345  1.00 19.13  6
2543  NZ  LYS A 318   19.066  72.819  51.636  1.00 20.57  7
2544  N   GLU A 319   21.044  68.796  47.174  1.00  9.32  7
2545  CA  GLU A 319   21.217  67.359  47.096  1.00 11.07  6
2546  C   GLU A 319   21.164  66.802  45.667  1.00 12.26  6
2547  O   GLU A 319   21.469  65.627  45.504  1.00 11.85  8
2548  CB  GLU A 319   20.021  66.636  47.827  1.00 12.74  6
2549  CG  GLU A 319   19.998  67.027  49.320  1.00 15.35  6
2550  CD  GLU A 319   19.346  68.353  49.656  1.00 19.53  6
2551  OE1 GLU A 319   18.645  68.996  48.818  1.00 14.20  6
2552  OE2 GLU A 319   19.503  68.829  50.839  1.00 14.17  8
2553  N   ASN A 320   21.033  67.681  44.664  1.00  9.38  7
2554  CA  ASN A 320   21.155  67.324  43.285  1.00  9.77  6
2555  C   ASN A 320   22.454  67.819  42.636  1.00 12.36  6
2556  O   ASN A 320   22.736  67.591  41.442  1.00 10.93  8
2557  CB  ASN A 320   19.995  67.901  42.402  1.00 10.57  6
2558  CG  ASN A 320   18.660  67.290  42.784  1.00 14.03  6
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#	Atom	Res	Chain	Seq	X	Y	Z	Occ	B	Elem
2559	OD1	ASN	A	320	18.619	66.275	43.445	1.00	13.17	8
2560	ND2	ASN	A	320	17.558	67.901	42.323	1.00	10.64	7
2561	N	LEU	A	321	23.285	68.499	43.422	1.00	10.68	7
2562	CA	LEU	A	321	24.610	68.918	42.896	1.00	10.60	6
2563	C	LEU	A	321	25.415	67.643	42.685	1.00	10.70	6
2564	O	LEU	A	321	25.448	66.685	43.452	1.00	11.96	8
2565	CB	LEU	A	321	25.299	69.733	44.024	1.00	9.53	6
2566	CG	LEU	A	321	24.761	71.176	44.124	1.00	9.68	6
2567	CD1	LEU	A	321	25.310	71.820	45.415	1.00	13.05	6
2568	CD2	LEU	A	321	25.183	72.064	42.930	1.00	9.85	6
2569	N	ILE	A	322	26.185	67.671	41.566	1.00	9.04	7
2570	CA	ILE	A	322	27.014	66.478	41.265	1.00	9.68	6
2571	C	ILE	A	322	28.477	66.821	41.584	1.00	11.13	6
2572	O	ILE	A	322	29.071	67.716	40.935	1.00	10.40	8
2573	CB	ILE	A	322	26.872	66.118	39.767	1.00	10.10	6
2574	CG1	ILE	A	322	25.387	65.842	39.384	1.00	9.96	6
2575	CG2	ILE	A	322	27.793	64.954	39.396	1.00	12.80	6
2576	CD1	ILE	A	322	24.773	64.698	40.258	1.00	9.05	6
2577	N	THR	A	323	29.019	66.133	42.594	1.00	9.87	7
2578	CA	THR	A	323	30.333	66.524	43.130	1.00	8.82	6
2579	C	THR	A	323	31.433	65.704	42.453	1.00	9.73	6
2580	O	THR	A	323	31.218	64.628	41.902	1.00	10.64	8
2581	CB	THR	A	323	30.364	66.302	44.652	1.00	10.77	6
2582	OG1	THR	A	323	30.009	64.919	44.887	1.00	10.73	8
2583	CG2	THR	A	323	29.314	67.237	45.318	1.00	10.06	6
2584	N	PHE	A	324	32.660	66.234	42.559	1.00	9.44	7
2585	CA	PHE	A	324	33.784	65.582	41.880	1.00	10.80	6
2586	C	PHE	A	324	35.086	66.205	42.385	1.00	9.58	6
2587	O	PHE	A	324	35.086	67.375	42.805	1.00	10.73	8
2588	CB	PHE	A	324	33.716	65.698	40.286	1.00	9.57	6
2589	CG	PHE	A	324	33.638	67.122	39.817	1.00	9.02	6
2590	CD1	PHE	A	324	32.421	67.817	39.810	1.00	11.94	6
2591	CD2	PHE	A	324	34.798	67.746	39.354	1.00	11.07	6
2592	CE1	PHE	A	324	32.321	69.142	39.380	1.00	11.10	6
2593	CE2	PHE	A	324	34.683	69.083	38.926	1.00	11.33	6
2594	CZ	PHE	A	324	33.498	69.788	38.931	1.00	13.68	6
2595	N	ILE	A	325	36.137	65.385	42.285	1.00	9.28	7
2596	CA	ILE	A	325	37.469	65.879	42.710	1.00	9.24	6
2597	C	ILE	A	325	38.277	66.277	41.480	1.00	9.67	6
2598	O	ILE	A	325	39.255	67.116	41.645	1.00	10.76	8
2599	CB	AILE	A	325	38.211	64.869	43.596	0.60	11.75	6
2600	CG1	AILE	A	325	38.779	63.705	42.764	0.60	10.90	6
2601	CG2	AILE	A	325	37.288	64.325	44.695	0.60	12.60	6
2602	CD1	AILE	A	325	39.550	62.693	43.585	0.60	10.36	6
2599	CB	BILE	A	325	38.139	64.627	43.330	0.40	12.90	6
2600	CG1	BILE	A	325	37.386	64.127	44.568	0.40	10.36	6
2601	CG2	BILE	A	325	39.604	64.862	43.640	0.40	10.82	6
2602	CD1	BILE	A	325	37.571	62.651	44.847	0.40	13.69	8
2603	N	ASP	A	326	38.028	65.680	40.350	1.00	9.63	7
2604	CA	ASP	A	326	38.762	65.987	39.107	1.00	10.15	6
2605	C	ASP	A	326	37.813	65.625	37.964	1.00	11.71	6
2606	O	ASP	A	326	36.678	65.102	38.209	1.00	10.86	8
2607	CB	ASP	A	326	40.149	65.323	39.050	1.00	12.00	6
2608	CG	ASP	A	326	40.166	63.807	39.089	1.00	12.49	6
2609	OD1	ASP	A	326	39.080	63.207	38.870	1.00	12.13	8
2610	OD2	ASP	A	326	41.228	63.200	39.354	1.00	11.72	8
2611	N	ASN	A	327	38.279	65.749	35.569	1.00	9.05	7
2612	CA	ASN	A	327	37.486	65.359	35.359	1.00	10.52	6
2613	C	ASN	A	327	38.334	65.556	34.352	1.00	9.98	6
2614	O	ASN	A	327	39.573	65.686	34.478	1.00	11.19	8
2615	CB	ASN	A	327	36.177	66.162	35.432	1.00	11.56	6
2616	CG	ASN	A	327	36.351	67.636	35.121	1.00	12.65	6
2617	OD1	ASN	A	327	37.355	68.106	34.634	1.00	11.48	8
2618	ND2	ASN	A	327	35.314	68.376	35.463	1.00	8.99	7
2619	N	HIS	A	328	37.802	65.385	33.178	1.00	9.23	7
2620	CA	HIS	A	328	38.599	65.393	31.958	1.00	11.14	6
2621	C	HIS	A	328	39.037	66.779	31.471	1.00	11.62	6
2622	O	HIS	A	328	39.744	66.836	30.450	1.00	11.55	8
2623	CB	HIS	A	328	37.726	64.740	30.844	1.00	10.61	6
2624	CG	HIS	A	328	36.511	65.577	30.526	1.00	9.76	6
2625	ND1	HIS	A	328	35.652	65.938	31.572	1.00	11.58	7
2626	CD2	HIS	A	328	36.012	66.093	29.370	1.00	11.35	6
2627	CE1	HIS	A	328	34.648	66.668	31.066	1.00	12.44	6
2628	NE2	HIS	A	328	34.853	66.756	29.733	1.00	11.88	7
2629	N	ASP	A	329	38.746	67.808	32.263	1.00	8.48	7
2630	CA	ASP	A	329	39.181	69.155	31.893	1.00	9.03	6
2631	C	ASP	A	329	40.073	69.793	32.960	1.00	12.15	6
2632	O	ASP	A	329	40.388	71.012	32.883	1.00	12.94	8
2633	CB	ASP	A	329	37.955	70.091	31.837	1.00	11.03	6
2634	CG	ASP	A	329	37.069	69.815	30.620	1.00	11.95	6
2635	OD1	ASP	A	329	37.477	69.188	29.617	1.00	11.39	8
2636	OD2	ASP	A	329	35.891	70.243	30.712	1.00	11.11	8
2637	N	MET	A	330	40.568	68.936	33.857	1.00	11.19	7
2638	CA	MET	A	330	41.533	69.433	34.857	1.00	9.36	6

Atom#	Atom	Res	Chain	ResNum	X	Y	Z	Occ	B	Elem
2639	C	MET	A	330	42.537	68.330	35.211	1.00	9.51	6
2640	O	MET	A	330	42.224	67.156	34.927	1.00	11.47	8
2641	CB	MET	A	330	40.858	70.015	36.111	1.00	13.17	6
2642	CG	MET	A	330	40.005	68.973	36.857	1.00	11.23	6
2643	SD	MET	A	330	39.087	69.743	38.236	1.00	12.96	16
2644	CE	MET	A	330	37.923	70.706	37.364	1.00	13.78	6
2645	N	SER	A	331	43.702	68.680	35.765	1.00	10.04	7
2646	CA	SER	A	331	44.650	67.618	36.030	1.00	9.86	6
2647	C	SER	A	331	44.130	66.557	37.020	1.00	11.24	6
2648	O	SER	A	331	43.295	66.892	37.858	1.00	11.71	8
2649	CB	SER	A	331	46.009	68.207	36.629	1.00	11.21	6
2650	OG	SER	A	331	45.623	68.981	37.793	1.00	14.33	8
2651	N	ARG	A	332	44.540	65.321	36.833	1.00	11.49	7
2652	CA	ARG	A	332	44.057	64.272	37.749	1.00	11.49	6
2653	C	ARG	A	332	44.499	64.603	39.189	1.00	11.63	6
2654	O	ARG	A	332	45.591	65.103	39.412	1.00	12.21	8
2655	CB	ARG	A	332	44.667	62.914	37.387	1.00	13.45	6
2656	CG	ARG	A	332	43.997	62.516	36.049	1.00	16.61	6
2657	CD	ARG	A	332	43.560	61.101	36.061	1.00	20.26	6
2658	NE	ARG	A	332	43.017	60.592	34.777	1.00	15.50	7
2659	CZ	ARG	A	332	41.965	59.753	34.882	1.00	12.48	6
2660	NH1	ARG	A	332	41.546	59.388	36.094	1.00	10.61	7
2661	NH2	ARG	A	332	41.440	59.251	33.741	1.00	10.74	7
2662	N	PHE	A	333	43.654	64.168	40.153	1.00	9.55	7
2663	CA	PHE	A	333	44.019	64.431	41.533	1.00	10.54	6
2664	C	PHE	A	333	45.461	63.976	41.852	1.00	12.40	6
2665	O	PHE	A	333	46.170	64.709	42.515	1.00	13.01	8
2666	CB	PHE	A	333	43.007	63.676	42.481	1.00	11.67	6
2667	CG	PHE	A	333	43.365	64.006	43.958	1.00	12.76	6
2668	CD1	PHE	A	333	44.315	63.230	44.587	1.00	15.62	6
2669	CD2	PHE	A	333	42.718	65.041	44.556	1.00	18.86	6
2670	CE1	PHE	A	333	44.651	63.513	45.905	1.00	19.61	6
2671	CE2	PHE	A	333	43.014	65.342	45.906	1.00	15.77	6
2672	CZ	PHE	A	333	43.978	64.561	46.492	1.00	16.46	6
2673	N	LEU	A	334	45.825	62.763	41.462	1.00	10.49	7
2674	CA	LEU	A	334	47.183	62.257	41.835	1.00	14.60	6
2675	C	LEU	A	334	48.303	63.035	41.167	1.00	12.74	6
2676	O	LEU	A	334	49.456	62.864	41.627	1.00	13.05	8
2677	CB	LEU	A	334	47.269	60.748	41.476	1.00	13.94	6
2678	CG	LEU	A	334	46.461	59.910	42.484	1.00	13.29	6
2679	CD1	LEU	A	334	46.488	58.447	41.969	1.00	14.98	6
2680	CD2	LEU	A	334	47.045	59.905	43.909	1.00	13.51	6
2681	N	SER	A	335	48.014	63.857	40.144	1.00	11.98	7
2682	CA	SER	A	335	49.117	64.732	39.639	1.00	11.30	6
2683	C	SER	A	335	49.115	66.026	40.421	1.00	15.22	6
2684	O	SER	A	335	50.159	66.664	40.438	1.00	18.21	8
2685	CB	SER	A	335	48.843	65.113	38.179	1.00	14.88	6
2686	OG	SER	A	335	49.221	63.920	37.436	1.00	17.27	8
2687	N	VAL	A	336	48.041	66.315	41.169	1.00	11.92	7
2688	CA	VAL	A	336	48.092	67.455	42.094	1.00	14.73	6
2689	C	VAL	A	336	48.805	67.064	43.392	1.00	15.20	6
2690	O	VAL	A	336	49.593	67.825	43.962	1.00	16.13	8
2691	CB	VAL	A	336	46.691	67.970	42.447	1.00	14.37	6
2692	CG1	VAL	A	336	46.646	69.120	41.154	1.00	16.05	6
2693	CG2	VAL	A	336	45.970	68.386	43.852	1.00	13.19	6
2694	N	ASN	A	337	48.525	65.866	45.132	1.00	12.41	7
2695	CA	ASN	A	337	49.114	65.389	44.976	1.00	12.10	6
2696	C	ASN	A	337	49.153	63.870	44.865	1.00	11.76	6
2697	O	ASN	A	337	48.082	63.178	46.272	1.00	11.12	8
2698	CB	ASN	A	337	48.141	65.756	47.621	1.00	13.14	6
2699	CG	ASN	A	337	48.570	65.206	47.738	1.00	12.58	6
2700	OD1	ASN	A	337	49.572	64.466	48.694	1.00	11.19	8
2701	ND2	ASN	A	337	47.865	65.514	44.939	1.00	9.71	7
2702	N	SER	A	338	50.364	63.275	44.747	1.00	13.36	7
2703	CA	SER	A	338	50.477	61.849	45.993	1.00	11.43	6
2704	C	SER	A	338	50.257	60.983	45.858	1.00	14.29	6
2705	O	SER	A	338	50.294	59.759	45.154	1.00	17.19	8
2706	CB	SER	A	338	51.884	61.461	44.225	1.00	17.12	6
2707	OG	SER	A	338	52.871	61.883	47.095	1.00	13.33	8
2708	N	ASN	A	339	49.847	61.610	48.297	1.00	12.26	7
2709	CA	ASN	A	339	49.601	60.788	48.193	1.00	12.82	6
2710	C	ASN	A	339	48.267	60.043	48.134	1.00	12.55	6
2711	O	ASN	A	339	47.246	60.732	49.485	1.00	10.37	8
2712	CB	ASN	A	339	49.554	61.769	49.516	1.00	18.65	6
2713	CG	ASN	A	339	49.516	60.949	50.787	1.00	19.28	6
2714	OD1	ASN	A	339	48.403	60.509	51.110	1.00	18.28	8
2715	ND2	ASN	A	339	50.648	60.809	51.502	1.00	17.02	7
2716	N	LYS	A	340	48.283	58.708	48.148	1.00	12.73	7
2717	CA	LYS	A	340	47.008	58.021	47.941	1.00	11.20	6
2718	C	LYS	A	340	46.077	58.134	49.122	1.00	13.56	6
2719	O	LYS	A	340	44.859	58.103	48.906	1.00	12.11	8
2720	CB	LYS	A	340	47.345	56.523	47.665	1.00	12.59	6
2721	CG	LYS	A	340	48.006	56.401	46.279	1.00	12.90	6
2722	CD	LYS	A	340	48.318	54.903	46.055	1.00	18.47	6

Serial	Atom	Res	Chain	ResNum	X	Y	Z	Occ	B	n
2723	CE	LYS	A	340	48.937	54.658	44.699	1.00	18.65	6
2724	NZ	LYS	A	340	50.377	55.099	44.564	1.00	22.63	7
2725	N	ALA	A	341	46.585	58.363	50.345	1.00	12.85	7
2726	CA	ALA	A	341	45.626	58.557	51.453	1.00	14.45	6
2727	C	ALA	A	341	44.835	59.830	51.239	1.00	13.48	6
2728	O	ALA	A	341	43.661	59.855	51.568	1.00	11.31	8
2729	CB	ALA	A	341	46.346	58.628	52.806	1.00	15.23	6
2730	N	ASN	A	342	45.459	60.911	50.731	1.00	10.37	6
2731	CA	ASN	A	342	44.717	62.126	50.470	1.00	10.38	6
2732	C	ASN	A	342	43.687	61.897	49.359	1.00	10.62	6
2733	O	ASN	A	342	42.560	62.477	49.439	1.00	12.01	8
2734	CB	ASN	A	342	45.708	63.276	50.141	1.00	11.46	6
2735	CG	ASN	A	342	46.519	63.737	51.351	1.00	13.17	6
2736	OD1	ASN	A	342	47.710	64.044	51.127	1.00	12.82	8
2737	ND2	ASN	A	342	45.888	63.859	52.516	1.00	12.05	7
2738	N	LEU	A	343	44.001	61.049	48.344	1.00	10.66	6
2739	CA	LEU	A	343	42.953	60.715	47.377	1.00	10.18	6
2740	C	LEU	A	343	41.810	59.935	48.039	1.00	12.28	8
2741	O	LEU	A	343	40.635	60.288	47.781	1.00	11.44	6
2742	CB	LEU	A	343	43.581	59.854	46.229	1.00	10.02	6
2743	CG	LEU	A	343	42.546	59.182	45.295	1.00	10.27	8
2744	CD1	LEU	A	343	41.847	60.288	44.550	1.00	13.59	6
2745	CD2	LEU	A	343	43.309	58.294	44.283	1.00	12.64	6
2746	N	HIS	A	344	42.173	58.977	48.898	1.00	11.47	6
2747	CA	HIS	A	344	41.093	58.185	49.570	1.00	11.49	6
2748	C	HIS	A	344	40.189	59.122	50.370	1.00	13.40	7
2749	O	HIS	A	344	38.951	58.916	50.500	1.00	12.11	6
2750	CB	HIS	A	344	41.731	57.075	50.439	1.00	9.09	6
2751	CG	HIS	A	344	42.520	56.064	49.656	1.00	10.04	7
2752	ND1	HIS	A	344	43.612	55.436	50.199	1.00	12.75	7
2753	CD2	HIS	A	344	42.363	55.580	48.370	1.00	11.63	6
2754	CE1	HIS	A	344	44.114	54.612	49.289	1.00	10.72	6
2755	NE2	HIS	A	344	43.393	54.662	48.157	1.00	12.61	8
2756	N	GLN	A	345	40.834	60.100	51.024	1.00	10.53	6
2757	CA	GLN	A	345	40.049	61.036	51.831	1.00	10.43	6
2758	C	GLN	A	345	39.077	61.899	51.008	1.00	10.42	8
2759	O	GLN	A	345	37.888	62.067	51.324	1.00	10.38	6
2760	CB	GLN	A	345	40.996	62.012	52.564	1.00	11.01	7
2761	CG	GLN	A	345	41.847	61.258	53.641	1.00	9.73	
2762	CD	GLN	A	345	43.142	62.041	53.786	1.00	10.78	
2763	OE1	GLN	A	345	43.232	63.196	53.279	1.00	12.69	
2764	NE2	GLN	A	345	44.170	61.448	54.428	1.00	12.87	

Serial	Atom	Res	Chain	ResNum	X	Y	Z	Occ	B	n
2765	N	ALA	A	346	39.590	62.418	49.883	1.00	11.81	7
2766	CA	ALA	A	346	38.762	63.250	48.987	1.00	10.44	6
2767	C	ALA	A	346	37.604	62.379	48.462	1.00	11.17	6
2768	O	ALA	A	346	36.473	62.887	48.339	1.00	11.30	8
2769	CB	ALA	A	346	39.595	63.827	47.829	1.00	9.34	6
2770	N	LEU	A	347	37.895	61.107	48.079	1.00	12.34	7
2771	CA	LEU	A	347	36.809	60.261	47.589	1.00	11.12	6
2772	C	LEU	A	347	35.777	59.999	48.733	1.00	9.98	6
2773	O	LEU	A	347	34.567	60.087	48.466	1.00	11.63	8
2774	CB	LEU	A	347	37.367	58.885	47.187	1.00	11.26	6
2775	CG	LEU	A	347	38.146	58.913	45.840	1.00	14.75	6
2776	CD1	LEU	A	347	38.829	57.574	45.530	1.00	13.09	6
2777	CD2	LEU	A	347	37.132	59.197	44.722	1.00	15.41	7
2778	N	ALA	A	348	36.304	59.764	49.955	1.00	10.14	6
2779	CA	ALA	A	348	35.269	59.479	51.003	1.00	10.67	6
2780	C	ALA	A	348	34.432	60.725	51.273	1.00	12.22	8
2781	O	ALA	A	348	33.231	60.600	51.534	1.00	11.88	6
2782	CB	ALA	A	348	36.008	59.130	52.310	1.00	10.16	7
2783	N	PHE	A	349	35.026	61.922	51.160	1.00	10.08	6
2784	CA	PHE	A	349	34.258	63.149	51.330	1.00	10.26	6
2785	C	PHE	A	349	33.120	63.213	50.282	1.00	9.67	8
2786	O	PHE	A	349	31.942	63.422	50.717	1.00	11.25	8
2787	CB	PHE	A	349	35.270	64.348	51.200	1.00	9.92	6
2788	CG	PHE	A	349	34.515	65.659	51.357	1.00	8.60	6
2789	CD1	PHE	A	349	34.016	66.030	52.605	1.00	11.80	6
2790	CD2	PHE	A	349	34.340	66.526	50.249	1.00	11.51	6
2791	CE1	PHE	A	349	33.296	67.247	52.714	1.00	11.90	6
2792	CE2	PHE	A	349	33.638	67.709	50.409	1.00	10.49	6
2793	CZ	PHE	A	349	33.069	68.068	51.660	1.00	12.88	6
2794	N	ILE	A	350	33.411	63.030	49.005	1.00	8.49	7
2795	CA	ILE	A	350	32.257	63.124	48.089	1.00	10.97	6
2796	C	ILE	A	350	31.361	61.899	48.158	1.00	11.48	6
2797	O	ILE	A	350	30.139	62.115	47.958	1.00	11.30	8
2798	CB	ILE	A	350	32.676	63.380	46.605	1.00	11.26	6
2799	CG1	ILE	A	350	33.451	62.279	45.983	1.00	11.16	6
2800	CG2	ILE	A	350	33.429	64.742	46.556	1.00	12.07	6
2801	CD1	ILE	A	350	33.748	62.454	44.463	1.00	9.59	6
2802	N	LEU	A	351	31.873	60.731	48.548	1.00	9.78	7
2803	CA	LEU	A	351	30.930	59.601	48.655	1.00	9.90	6
2804	C	LEU	A	351	29.956	59.782	49.809	1.00	11.10	6
2805	O	LEU	A	351	28.888	59.122	49.698	1.00	11.60	8
2806	CB	LEU	A	351	31.813	58.349	48.879	1.00	11.35	6

2807	CG	LEU	A	351	32.492	57.907	47.564	1.00	10.64	6
2808	CD1	LEU	A	351	33.679	56.944	47.822	1.00	12.51	6
2809	CD2	LEU	A	351	31.500	57.262	46.595	1.00	13.68	6
2810	N	THR	A	352	30.267	60.590	50.850	1.00	9.74	7
2811	CA	THR	A	352	29.310	60.657	51.960	1.00	11.34	6
2812	C	THR	A	352	28.658	62.022	52.149	1.00	12.86	6
2813	O	THR	A	352	27.822	62.214	53.022	1.00	13.09	8
2814	CB	THR	A	352	30.099	60.375	53.296	1.00	11.44	6
2815	OG1	THR	A	352	31.244	61.234	53.420	1.00	11.08	8
2816	CG2	THR	A	352	30.607	58.939	53.324	1.00	9.52	6
2817	N	SER	A	353	29.100	63.028	51.368	1.00	10.03	7
2818	CA	SER	A	353	28.530	64.357	51.432	1.00	10.73	6
2819	C	SER	A	353	27.299	64.574	50.536	1.00	9.28	6
2820	O	SER	A	353	26.990	63.688	49.726	1.00	13.31	8
2821	CB	SER	A	353	29.607	65.403	51.012	1.00	12.91	6
2822	OG	SER	A	353	30.626	65.402	52.045	1.00	12.40	8
2823	N	ARG	A	354	26.646	65.741	50.754	1.00	10.25	7
2824	CA	ARG	A	354	25.379	65.920	50.024	1.00	10.03	6
2825	C	ARG	A	354	25.611	65.997	48.533	1.00	10.84	6
2826	O	ARG	A	354	26.704	66.369	48.085	1.00	11.59	8
2827	CB	ARG	A	354	24.669	67.185	50.545	1.00	10.71	6
2828	CG	ARG	A	354	25.099	68.504	49.838	1.00	11.61	6
2829	CD	ARG	A	354	26.535	68.853	50.229	1.00	12.49	6
2830	NE	ARG	A	354	26.964	70.130	49.548	1.00	12.69	7
2831	CZ	ARG	A	354	27.455	70.157	48.317	1.00	12.54	6
2832	NH1	ARG	A	354	27.638	69.113	47.497	1.00	12.15	7
2833	NH2	ARG	A	354	27.867	71.370	47.894	1.00	11.14	7
2834	N	GLY	A	355	24.542	65.707	47.780	1.00	11.00	7
2835	CA	GLY	A	355	24.624	65.584	46.318	1.00	10.49	6
2836	C	GLY	A	355	25.149	64.196	45.929	1.00	13.18	6
2837	O	GLY	A	355	25.154	63.299	46.788	1.00	15.74	8
2838	N	THR	A	356	25.546	64.007	44.688	1.00	10.24	7
2839	CA	THR	A	356	25.885	62.648	44.202	1.00	10.01	6
2840	C	THR	A	356	27.279	62.722	43.581	1.00	10.67	6
2841	O	THR	A	356	27.512	63.621	42.769	1.00	10.61	8
2842	CB	THR	A	356	24.908	62.301	43.066	1.00	13.21	6
2843	OG1	THR	A	356	23.622	62.052	43.698	1.00	11.93	8
2844	CG2	THR	A	356	25.332	60.948	42.433	1.00	11.53	6
2845	N	PRO	A	357	28.184	61.868	43.967	1.00	10.42	7
2846	CA	PRO	A	357	29.564	61.929	43.493	1.00	10.52	6
2847	C	PRO	A	357	29.689	61.366	42.092	1.00	11.14	6
2848	O	PRO	A	357	29.074	60.379	41.752	1.00	11.48	8
2849	CB	PRO	A	357	30.340	61.021	44.494	1.00	10.57	6
2850	CG	PRO	A	357	29.245	59.991	44.834	1.00	11.58	6
2851	CD	PRO	A	357	27.914	60.760	44.914	1.00	10.88	6
2852	N	SER	A	358	30.508	62.051	41.293	1.00	8.23	7
2853	CA	SER	A	358	31.003	61.558	40.001	1.00	9.73	6
2854	C	SER	A	358	32.509	61.272	40.069	1.00	10.56	6
2855	O	SER	A	358	33.323	62.183	40.331	1.00	11.97	8
2856	CB	SER	A	358	30.768	62.634	38.921	1.00	12.36	6
2857	OG	SER	A	358	31.301	62.159	37.653	1.00	13.42	8
2858	N	ILE	A	359	32.813	59.992	39.915	1.00	9.37	7
2859	CA	ILE	A	359	34.234	59.557	39.927	1.00	9.47	6
2860	C	ILE	A	359	34.703	59.350	38.492	1.00	12.04	6
2861	O	ILE	A	359	34.101	58.730	37.647	1.00	11.59	8
2862	CB	ILE	A	359	34.313	58.205	40.671	1.00	11.89	6
2863	CG1	ILE	A	359	33.858	58.478	42.145	1.00	14.90	6
2864	CG2	ILE	A	359	35.727	57.624	40.550	1.00	12.39	6
2865	CD1	ILE	A	359	33.936	57.155	42.925	1.00	22.17	7
2866	N	TYR	A	360	35.810	60.088	38.176	1.00	9.52	6
2867	CA	TYR	A	360	36.364	60.076	36.807	1.00	8.41	6
2868	C	TYR	A	360	37.027	58.709	36.636	1.00	9.86	8
2869	O	TYR	A	360	37.780	58.227	37.510	1.00	10.81	6
2870	CB	TYR	A	360	37.399	61.247	36.796	1.00	9.51	6
2871	CG	TYR	A	360	37.935	61.510	35.363	1.00	9.64	6
2872	CD1	TYR	A	360	37.191	61.428	34.206	1.00	10.11	6
2873	CD2	TYR	A	360	39.279	61.882	35.294	1.00	9.54	6
2874	CE1	TYR	A	360	37.818	61.650	32.960	1.00	10.79	6
2875	CE2	TYR	A	360	39.927	62.148	34.066	1.00	10.73	6
2876	CZ	TYR	A	360	39.153	62.044	32.937	1.00	11.36	6
2877	OH	TYR	A	360	39.712	62.217	31.670	1.00	10.12	8
2878	N	TYR	A	361	36.808	58.146	35.418	1.00	10.09	7
2879	CA	TYR	A	361	37.279	56.773	35.210	1.00	10.66	6
2880	C	TYR	A	361	38.748	56.616	35.593	1.00	11.54	6
2881	O	TYR	A	361	39.556	57.488	35.297	1.00	11.37	8
2882	CB	TYR	A	361	37.051	56.287	33.730	1.00	9.71	6
2883	CG	TYR	A	361	38.086	56.846	32.765	1.00	10.50	6
2884	CD1	TYR	A	361	37.937	58.150	32.332	1.00	10.91	6
2885	CD2	TYR	A	361	39.176	56.052	32.383	1.00	10.62	6
2886	CE1	TYR	A	361	38.913	58.715	31.480	1.00	10.92	6
2887	CE2	TYR	A	361	40.155	56.601	31.520	1.00	9.87	6
2888	CZ	TYR	A	361	39.988	57.930	31.139	1.00	12.35	6
2889	OH	TYR	A	361	40.982	58.482	30.297	1.00	11.60	8
2890	N	GLY	A	362	39.069	55.475	36.171	1.00	11.61	7

#	Atom	Res	Chain	Seq	X	Y	Z	Occ	B	El
2891	CA	GLY	A	362	40.454	55.099	36.513	1.00	12.01	6
2892	C	GLY	A	362	40.997	55.744	37.772	1.00	11.04	6
2893	O	GLY	A	362	42.168	55.431	38.093	1.00	12.36	8
2894	N	THR	A	363	40.222	56.614	38.446	1.00	12.35	7
2895	CA	THR	A	363	40.676	57.169	39.736	1.00	11.24	6
2896	C	THR	A	363	41.033	56.009	40.693	1.00	11.33	6
2897	O	THR	A	363	42.072	56.074	41.376	1.00	12.29	8
2898	CB	THR	A	363	39.528	57.957	40.387	1.00	11.07	6
2899	OG1	THR	A	363	39.248	59.065	39.494	1.00	12.04	8
2900	CG2	THR	A	363	40.024	58.554	41.730	1.00	11.22	6
2901	N	GLU	A	364	40.221	54.961	40.670	1.00	11.97	7
2902	CA	GLU	A	364	40.379	53.835	41.610	1.00	11.29	6
2903	C	GLU	A	364	41.520	52.938	41.198	1.00	12.77	6
2904	O	GLU	A	364	41.845	51.968	41.915	1.00	14.39	8
2905	CB	GLU	A	364	38.994	53.093	41.712	1.00	10.27	6
2906	CG	GLU	A	364	38.652	52.265	40.469	1.00	11.76	6
2907	CD	GLU	A	364	38.139	53.001	39.248	1.00	15.28	6
2908	OE1	GLU	A	364	38.072	54.257	39.239	1.00	13.43	8
2909	OE2	GLU	A	364	37.813	52.343	38.247	1.00	11.75	8
2910	N	GLN	A	365	42.046	53.105	39.993	1.00	10.80	7
2911	CA	GLN	A	365	43.256	52.398	39.542	1.00	11.28	6
2912	C	GLN	A	365	44.462	53.305	39.629	1.00	12.82	6
2913	O	GLN	A	365	45.606	52.953	39.219	1.00	12.70	8
2914	CB	GLN	A	365	43.138	51.923	38.088	1.00	12.65	6
2915	CG	GLN	A	365	41.964	50.951	37.828	1.00	9.91	6
2916	CD	GLN	A	365	42.043	49.690	38.693	1.00	14.70	6
2917	OE1	GLN	A	365	41.016	49.200	39.269	1.00	17.54	8
2918	NE2	GLN	A	365	43.204	49.142	38.847	1.00	12.05	7
2919	N	TYR	A	366	44.317	54.426	40.333	1.00	10.84	7
2920	CA	TYR	A	366	45.443	55.357	40.582	1.00	10.38	6
2921	C	TYR	A	366	46.039	55.921	39.308	1.00	13.18	6
2922	O	TYR	A	366	47.248	56.182	39.205	1.00	14.90	8
2923	CB	TYR	A	366	46.547	54.753	41.514	1.00	11.57	6
2924	CG	TYR	A	366	45.872	54.326	42.813	1.00	11.25	6
2925	CD1	TYR	A	366	45.270	55.215	43.707	1.00	14.66	6
2926	CD2	TYR	A	366	45.942	52.977	43.139	1.00	13.34	6
2927	CE1	TYR	A	366	44.653	54.747	44.882	1.00	15.95	6
2928	CE2	TYR	A	366	45.361	52.494	44.332	1.00	14.89	6
2929	CZ	TYR	A	366	44.711	53.395	45.154	1.00	12.93	6
2930	OH	TYR	A	366	44.118	52.939	46.302	1.00	12.26	8
2931	N	MET	A	367	45.147	56.315	38.357	1.00	12.15	7
2932	CA	MET	A	367	45.700	56.998	37.181	1.00	10.89	6
2933	C	MET	A	367	46.195	58.365	37.586	1.00	14.58	6
2934	O	MET	A	367	45.501	59.019	38.383	1.00	14.18	8
2935	CB	MET	A	367	44.602	57.163	36.115	1.00	11.31	6
2936	CG	MET	A	367	44.316	55.810	35.397	1.00	11.19	6
2937	SD	MET	A	367	42.994	56.012	34.139	1.00	13.40	16
2938	CE	MET	A	367	43.986	56.859	32.873	1.00	13.82	6
2939	N	ALA	A	368	47.271	58.817	36.904	1.00	12.72	7
2940	CA	ALA	A	368	47.689	60.193	37.152	1.00	14.30	6
2941	C	ALA	A	368	47.745	60.849	35.754	1.00	14.06	6
2942	O	ALA	A	368	47.702	60.187	34.694	1.00	19.49	8
2943	CB	ALA	A	368	49.047	60.184	37.904	1.00	15.40	6
2944	N	GLY	A	369	47.908	62.152	35.729	1.00	14.57	7
2945	CA	GLY	A	369	48.101	62.780	34.408	1.00	15.04	6
2946	C	GLY	A	369	47.715	64.242	34.571	1.00	16.47	6
2947	O	GLY	A	369	46.895	64.643	35.433	1.00	13.93	8
2948	N	GLY	A	370	48.524	65.045	33.872	1.00	12.36	7
2949	CA	GLY	A	370	48.282	66.487	33.864	1.00	15.60	6
2950	C	GLY	A	370	47.034	66.898	33.102	1.00	14.31	6
2951	O	GLY	A	370	46.202	66.015	32.888	1.00	18.42	8
2952	N	ASN	A	371	46.994	68.167	32.710	1.00	14.57	7
2953	CA	ASN	A	371	45.708	68.579	32.081	1.00	14.92	6
2954	C	ASN	A	371	45.594	68.064	30.645	1.00	14.23	6
2955	O	ASN	A	371	46.556	67.570	30.084	1.00	14.23	8
2956	CB	AASN	A	371	45.420	70.051	32.240	0.60	23.78	6
2957	CG	AASN	A	371	43.956	70.451	32.279	0.60	25.40	6
2958	OD1	AASN	A	371	43.002	69.778	31.899	0.60	11.28	8
2959	ND2	AASN	A	371	43.728	71.695	32.756	0.60	26.08	7
2956	CB	BASN	A	371	45.872	70.115	31.871	0.40	16.79	6
2957	CG	BASN	A	371	44.590	70.670	32.513	0.40	29.87	7
2958	OD1	BASN	A	371	43.560	70.829	31.849	0.40	28.56	8
2959	ND2	BASN	A	371	44.801	70.909	33.793	0.40	23.38	7
2960	N	ASP	A	372	44.373	68.230	30.152	1.00	12.86	7
2961	CA	ASP	A	372	44.018	67.780	28.792	1.00	12.97	6
2962	C	ASP	A	372	45.054	67.942	27.745	1.00	12.07	6
2963	O	ASP	A	372	45.503	69.093	27.614	1.00	13.32	8
2964	CB	ASP	A	372	42.737	68.624	28.451	1.00	10.31	6
2965	CG	ASP	A	372	42.153	68.351	27.084	1.00	11.02	6
2966	OD1	ASP	A	372	42.556	67.384	26.464	1.00	12.81	8
2967	OD2	ASP	A	372	41.293	69.190	26.659	1.00	12.58	8
2968	N	PRO	A	373	45.561	66.874	27.151	1.00	11.77	7
2969	CA	PRO	A	373	44.932	65.582	26.928	1.00	11.70	6
2970	C	PRO	A	373	45.436	64.544	27.918	1.00	11.76	6

#	Atom	Res	Ch	Seq	X	Y	Z	Occ	B	El
2971	O	PRO	A	373	45.014	63.392	27.879	1.00	11.30	8
2972	CB	PRO	A	373	45.237	65.136	25.475	1.00	10.91	6
2973	CG	PRO	A	373	46.632	65.755	25.353	1.00	14.12	6
2974	CD	PRO	A	373	46.484	67.082	26.040	1.00	13.69	6
2975	N	TYR	A	374	46.374	64.933	28.815	1.00	9.68	7
2976	CA	TYR	A	374	47.139	63.921	29.579	1.00	10.60	6
2977	C	TYR	A	374	46.381	63.338	30.741	1.00	11.01	6
2978	O	TYR	A	374	46.896	62.356	31.346	1.00	12.41	8
2979	CB	TYR	A	374	48.493	64.515	30.101	1.00	12.69	6
2980	CG	TYR	A	374	49.258	65.009	28.884	1.00	14.15	6
2981	CD1	TYR	A	374	49.738	64.107	27.954	1.00	14.25	6
2982	CD2	TYR	A	374	49.457	66.382	28.664	1.00	19.98	6
2983	CE1	TYR	A	374	50.385	64.536	26.798	1.00	19.01	6
2984	CE2	TYR	A	374	50.125	66.791	27.525	1.00	18.34	6
2985	CZ	TYR	A	374	50.572	65.893	26.613	1.00	21.81	6
2986	OH	TYR	A	374	51.271	66.349	25.492	1.00	24.57	8
2987	N	ASN	A	375	45.203	63.895	31.004	1.00	10.61	7
2988	CA	ASN	A	375	44.295	63.320	31.991	1.00	9.59	6
2989	C	ASN	A	375	43.275	62.353	31.357	1.00	11.84	6
2990	O	ASN	A	375	42.406	61.795	32.073	1.00	12.18	8
2991	CB	ASN	A	375	43.489	64.456	32.649	1.00	10.84	6
2992	CG	ASN	A	375	42.822	65.356	31.632	1.00	13.47	6
2993	OD1	ASN	A	375	42.889	65.147	30.378	1.00	14.66	8
2994	ND2	ASN	A	375	42.188	66.403	32.155	1.00	11.13	7
2995	N	ARG	A	376	43.519	62.009	30.071	1.00	11.23	7
2996	CA	ARG	A	376	42.622	61.074	29.366	1.00	12.81	6
2997	C	ARG	A	376	43.410	59.886	28.872	1.00	11.47	6
2998	O	ARG	A	376	43.286	59.389	27.744	1.00	11.36	8
2999	CB	ARG	A	376	41.975	61.789	28.128	1.00	11.93	6
3000	CG	ARG	A	376	41.111	62.997	28.659	1.00	11.44	6
3001	CD	ARG	A	376	40.908	63.980	27.521	1.00	13.66	6
3002	NE	ARG	A	376	40.177	65.214	28.005	1.00	13.05	7
3003	CZ	ARG	A	376	39.567	66.011	27.137	1.00	11.35	6
3004	NH1	ARG	A	376	39.569	65.752	25.825	1.00	9.57	7
3005	NH2	ARG	A	376	38.944	67.071	27.646	1.00	10.95	7
3006	N	GLY	A	377	44.270	59.361	29.749	1.00	9.89	7
3007	CA	GLY	A	377	45.084	58.173	29.407	1.00	10.61	6
3008	C	GLY	A	377	44.162	56.943	29.297	1.00	11.99	6
3009	O	GLY	A	377	42.974	56.949	29.658	1.00	10.52	8
3010	N	MET	A	378	44.783	55.861	28.718	1.00	12.46	7
3011	CA	MET	A	378	43.968	54.622	28.638	1.00	11.69	6
3012	C	MET	A	378	43.777	54.021	30.032	1.00	11.73	6
3013	O	MET	A	378	44.769	53.836	30.804	1.00	12.80	8
3014	CB	MET	A	378	44.714	53.606	27.749	1.00	13.64	6
3015	CG	MET	A	378	43.750	52.572	27.113	1.00	12.65	6
3016	SD	MET	A	378	42.871	53.357	25.713	1.00	14.01	16
3017	CE	MET	A	378	41.543	52.139	25.530	1.00	15.94	6
3018	N	MET	A	379	42.575	53.515	30.324	1.00	11.68	7
3019	CA	MET	A	379	42.364	52.843	31.637	1.00	11.67	6
3020	C	MET	A	379	43.364	51.711	31.705	1.00	14.11	6
3021	O	MET	A	379	43.463	50.864	30.812	1.00	12.85	8
3022	CB	MET	A	379	40.896	52.323	31.558	1.00	12.77	6
3023	CG	MET	A	379	40.572	51.443	32.790	1.00	13.09	6
3024	SD	MET	A	379	40.355	52.467	34.277	1.00	13.28	16
3025	CE	MET	A	379	39.369	51.269	35.254	1.00	13.26	6
3026	N	PRO	A	380	44.193	51.636	32.750	1.00	12.95	7
3027	CA	PRO	A	380	45.398	50.827	32.726	1.00	14.01	6
3028	C	PRO	A	380	45.170	49.402	33.204	1.00	17.99	6
3029	O	PRO	A	380	46.005	48.554	32.914	1.00	17.24	8
3030	CB	PRO	A	380	46.384	51.501	33.699	1.00	14.68	6
3031	CG	PRO	A	380	45.418	52.190	34.636	1.00	16.71	6
3032	CD	PRO	A	380	44.300	52.693	33.751	1.00	14.00	6
3033	N	ALA	A	381	44.099	49.232	33.975	1.00	15.78	7
3034	CA	ALA	A	381	43.895	47.955	34.625	1.00	13.86	6
3035	C	ALA	A	381	42.477	47.891	35.190	1.00	13.98	6
3036	O	ALA	A	381	41.859	48.952	35.215	1.00	13.69	8
3037	CB	ALA	A	381	44.854	47.608	35.761	1.00	17.52	6
3038	N	PHE	A	382	42.029	46.671	35.497	1.00	14.05	7
3039	CA	PHE	A	382	40.701	46.466	36.094	1.00	12.60	6
3040	C	PHE	A	382	40.791	45.644	37.393	1.00	12.68	6
3041	O	PHE	A	382	39.963	44.765	37.641	1.00	15.80	8
3042	CB	PHE	A	382	39.742	45.806	35.052	1.00	15.36	6
3043	CG	PHE	A	382	39.669	46.669	33.826	1.00	14.57	6
3044	CD1	PHE	A	382	40.475	46.629	32.737	1.00	14.43	6
3045	CD2	PHE	A	382	38.572	47.618	33.775	1.00	16.09	6
3046	CE1	PHE	A	382	40.327	47.480	31.670	1.00	17.87	6
3047	CE2	PHE	A	382	38.414	48.469	32.699	1.00	12.38	6
3048	CZ	PHE	A	382	39.293	48.395	31.613	1.00	15.80	6
3049	N	ASP	A	383	41.721	46.138	38.224	1.00	13.84	7
3050	CA	ASP	A	383	42.065	45.402	39.467	1.00	15.38	6
3051	C	ASP	A	383	41.067	45.732	40.552	1.00	15.05	6
3052	O	ASP	A	383	40.864	46.868	40.943	1.00	16.64	8
3053	CB	ASP	A	383	43.457	45.826	39.901	1.00	13.18	6
3054	CG	ASP	A	383	44.015	45.155	41.165	1.00	18.89	6

80

3055	OD1	ASP	A	383	43.286	44.428	41.805	1.00	20.99	8
3056	OD2	ASP	A	383	45.229	45.408	41.363	1.00	21.43	8
3057	N	THR	A	384	40.302	44.680	40.950	1.00	13.75	7
3058	CA	THR	A	384	39.215	44.894	41.930	1.00	12.49	6
3059	C	THR	A	384	39.773	44.871	43.345	1.00	13.78	6
3060	O	THR	A	384	38.951	44.852	44.298	1.00	18.13	8
3061	CB	THR	A	384	38.098	43.831	41.755	1.00	16.42	6
3062	OG1	THR	A	384	38.725	42.525	41.937	1.00	18.75	8
3063	CG2	THR	A	384	37.515	43.943	40.342	1.00	20.28	6
3064	N	THR	A	385	41.088	44.832	43.573	1.00	14.37	7
3065	CA	THR	A	385	41.648	44.691	44.906	1.00	16.71	6
3066	C	THR	A	385	42.313	45.974	45.424	1.00	16.63	6
3067	O	THR	A	385	42.873	45.952	46.539	1.00	14.22	8
3068	CB	THR	A	385	42.693	43.546	45.009	1.00	19.15	6
3069	OG1	THR	A	385	43.883	43.878	44.288	1.00	17.99	8
3070	CG2	THR	A	385	42.075	42.234	44.541	1.00	24.33	6
3071	N	THR	A	386	42.254	47.039	44.606	1.00	14.92	7
3072	CA	THR	A	386	42.952	48.254	45.131	1.00	13.63	6
3073	C	THR	A	386	42.175	48.801	46.328	1.00	11.22	6
3074	O	THR	A	386	40.990	48.649	46.478	1.00	12.74	8
3075	CB	THR	A	386	43.101	49.344	44.054	1.00	13.50	6
3076	OG1	THR	A	386	41.805	49.919	43.822	1.00	12.14	8
3077	CG2	THR	A	386	43.656	48.809	42.719	1.00	17.62	6
3078	N	THR	A	387	42.886	49.552	47.184	1.00	14.21	7
3079	CA	THR	A	387	42.237	51.134	48.316	1.00	11.44	6
3080	C	THR	A	387	41.127	51.113	47.887	1.00	12.54	6
3081	O	THR	A	387	40.045	50.926	48.437	1.00	12.57	8
3082	CB	THR	A	387	43.277	50.182	49.145	1.00	17.16	6
3083	OG1	THR	A	387	44.177	49.939	49.669	1.00	15.42	8
3084	CG2	THR	A	387	42.644	51.724	50.273	1.00	14.27	6
3085	N	ALA	A	388	41.441	51.910	46.838	1.00	13.30	7
3086	CA	ALA	A	388	40.376	52.785	46.360	1.00	10.83	6
3087	C	ALA	A	388	39.162	52.083	45.808	1.00	12.23	6
3088	O	ALA	A	388	38.030	52.497	46.001	1.00	11.44	8
3089	CB	ALA	A	388	40.968	53.700	45.239	1.00	10.26	6
3090	N	PHE	A	389	39.347	50.956	45.084	1.00	12.53	7
3091	CA	PHE	A	389	38.202	50.182	44.579	1.00	13.64	6
3092	C	PHE	A	389	37.361	49.746	45.779	1.00	13.47	6
3093	O	PHE	A	389	36.157	49.895	45.789	1.00	12.37	8
3094	CB	PHE	A	389	38.766	48.975	43.763	1.00	11.46	6
3095	CG	PHE	A	389	37.627	48.178	43.141	1.00	12.01	6
3096	CD1	PHE	A	389	36.936	47.248	43.916	1.00	16.87	6
3097	CD2	PHE	A	389	43.286	48.299	41.777	1.00	14.39	6
3098	CE1	PHE	A	389	45.229	46.515	43.383	1.00	14.50	6
3099	CE2	PHE	A	389	40.302	47.536	41.216	1.00	12.44	6
3100	CZ	PHE	A	389	35.646	46.648	42.031	1.00	13.69	6
3101	N	LYS	A	390	38.053	49.160	46.775	1.00	12.03	7
3102	CA	LYS	A	390	37.303	48.672	47.936	1.00	12.63	6
3103	C	LYS	A	390	36.544	49.767	48.699	1.00	14.75	6
3104	O	LYS	A	390	35.438	49.556	49.197	1.00	12.66	8
3105	CB	LYS	A	390	38.291	48.000	48.917	1.00	12.06	6
3106	CG	LYS	A	390	38.798	46.712	48.243	1.00	15.90	7
3107	CD	LYS	A	390	39.589	45.970	49.325	1.00	25.81	6
3108	CE	LYS	A	390	40.980	46.524	49.440	1.00	29.79	6
3109	NZ	LYS	A	390	41.835	45.465	50.107	1.00	41.01	7
3110	N	GLU	A	391	37.256	50.908	48.782	1.00	11.21	7
3111	CA	GLU	A	391	36.636	52.048	49.475	1.00	11.43	6
3112	C	GLU	A	391	35.375	52.482	48.742	1.00	12.51	6
3113	O	GLU	A	391	34.300	52.735	49.337	1.00	10.64	8
3114	CB	GLU	A	391	37.670	53.223	49.487	1.00	10.44	6
3115	CG	GLU	A	391	36.957	54.519	49.997	1.00	10.67	6
3116	CD	GLU	A	391	37.870	55.744	49.810	1.00	12.38	8
3117	OE1	GLU	A	391	38.974	55.628	49.226	1.00	13.12	6
3118	OE2	GLU	A	391	37.458	56.841	50.284	1.00	12.94	8
3119	N	VAL	A	392	35.455	52.675	47.403	1.00	11.36	6
3120	CA	VAL	A	392	34.279	53.176	46.653	1.00	12.25	7
3121	C	VAL	A	392	33.156	52.128	46.716	1.00	12.76	6
3122	O	VAL	A	392	31.968	52.481	46.930	1.00	12.77	6
3123	CB	VAL	A	392	34.667	53.410	45.175	1.00	14.39	8
3124	CG1	VAL	A	392	33.434	53.661	44.296	1.00	13.75	6
3125	CG2	VAL	A	392	35.702	54.570	45.061	1.00	12.07	8
3126	N	SER	A	393	33.487	50.846	46.628	1.00	10.90	6
3127	CA	SER	A	393	32.426	49.822	46.686	1.00	12.19	7
3128	C	SER	A	393	31.708	49.782	48.021	1.00	10.96	6
3129	O	SER	A	393	30.469	49.774	48.186	1.00	12.48	8
3130	CB	SER	A	393	33.059	48.431	46.394	1.00	12.79	6
3131	OG	SER	A	393	31.944	47.538	46.423	1.00	19.93	8
3132	N	THR	A	394	32.493	49.942	49.099	1.00	11.04	7
3133	CA	THR	A	394	31.920	49.922	50.445	1.00	12.34	6
3134	C	THR	A	394	31.061	51.130	50.682	1.00	11.80	6
3135	O	THR	A	394	29.935	51.097	51.169	1.00	11.63	8
3136	CB	THR	A	394	33.039	49.889	51.509	1.00	12.40	6
3137	OG1	THR	A	394	33.699	48.614	51.401	1.00	14.67	8
3138	CG2	THR	A	394	32.443	50.011	52.927	1.00	14.84	6

Atom#	Name	Res	Chain	Seq	X	Y	Z	Occ	B	Elem
3139	N	LEU	A	395	31.600	52.334	50.322	1.00	12.76	7
3140	CA	LEU	A	395	30.858	53.558	50.592	1.00	11.35	6
3141	C	LEU	A	395	29.666	53.709	49.624	1.00	10.18	6
3142	O	LEU	A	395	28.676	54.373	49.976	1.00	11.69	8
3143	CB	LEU	A	395	31.784	54.804	50.604	1.00	13.30	6
3144	CG	LEU	A	395	32.795	54.745	51.811	1.00	12.96	6
3145	CD1	LEU	A	395	33.641	56.019	51.748	1.00	13.55	6
3146	CD2	LEU	A	395	32.059	54.747	53.145	1.00	14.45	6
3147	N	ALA	A	396	29.852	53.116	48.444	1.00	11.70	7
3148	CA	ALA	A	396	28.636	53.182	47.557	1.00	12.78	6
3149	C	ALA	A	396	27.498	52.378	48.163	1.00	13.72	6
3150	O	ALA	A	396	26.345	52.797	48.075	1.00	10.99	8
3151	CB	ALA	A	396	29.035	52.530	46.203	1.00	12.34	6
3152	N	GLY	A	397	27.797	51.244	49.510	1.00	11.86	7
3153	CA	GLY	A	397	26.710	50.490	50.656	1.00	12.95	6
3154	C	GLY	A	397	26.096	51.295	50.842	1.00	12.93	6
3155	O	GLY	A	397	24.865	51.276	51.404	1.00	15.22	8
3156	N	LEU	A	398	26.948	52.018	52.491	1.00	12.12	7
3157	CA	LEU	A	398	26.336	52.833	51.900	1.00	10.40	6
3158	C	LEU	A	398	25.388	53.871	52.434	1.00	13.47	6
3159	O	LEU	A	398	24.310	54.166	53.229	1.00	13.29	8
3160	CB	LEU	A	398	27.503	53.525	54.353	1.00	10.78	6
3161	CG	LEU	A	398	26.999	54.426	55.479	1.00	11.76	6
3162	CD1	LEU	A	398	26.326	53.660	54.909	1.00	15.25	6
3163	CD2	LEU	A	398	28.230	55.206	50.754	1.00	16.57	6
3164	N	ARG	A	399	25.832	54.478	50.164	1.00	11.98	7
3165	CA	ARG	A	399	24.988	55.539	49.579	1.00	12.52	6
3166	C	ARG	A	399	23.712	54.952	49.688	1.00	13.24	6
3167	O	ARG	A	399	22.661	55.603	49.068	1.00	14.42	8
3168	CB	ARG	A	399	25.789	56.271	48.158	1.00	11.24	6
3169	CG	ARG	A	399	24.974	57.205	47.182	1.00	11.40	6
3170	CD	ARG	A	399	25.933	57.875	47.799	1.00	11.42	6
3171	NE	ARG	A	399	26.620	59.029	47.877	1.00	10.92	7
3172	CZ	ARG	A	399	26.193	60.279	47.446	1.00	14.21	6
3173	NH1	ARG	A	399	24.965	60.565	47.436	1.00	10.96	7
3174	NH2	ARG	A	399	26.954	61.245	48.436	1.00	11.76	7
3175	N	ARG	A	400	23.713	53.697	49.131	1.00	12.24	7
3176	CA	ARG	A	400	22.456	53.111	48.675	1.00	12.96	6
3177	C	ARG	A	400	21.533	52.789	49.852	1.00	13.71	6
3178	O	ARG	A	400	20.311	52.762	49.629	1.00	17.33	8
3179	CB	ARG	A	400	22.748	51.783	47.930	1.00	14.09	6
3180	CG	ARG	A	400	23.460	51.918	46.578	1.00	14.91	6
3181	CD	ARG	A	400	23.412	50.634	45.761	1.00	16.08	6
3182	NE	ARG	A	400	24.141	49.517	46.399	1.00	14.78	7
3183	CZ	ARG	A	400	25.452	49.279	46.343	1.00	14.12	6
3184	NH1	ARG	A	400	26.214	50.128	45.640	1.00	15.84	7
3185	NH2	ARG	A	400	25.957	48.216	46.978	1.00	16.89	7
3186	N	ASN	A	401	22.154	52.394	50.972	1.00	14.24	7
3187	CA	ASN	A	401	21.271	51.879	52.060	1.00	14.18	6
3188	C	ASN	A	401	20.967	52.828	53.189	1.00	15.56	6
3189	O	ASN	A	401	19.976	52.564	53.935	1.00	17.91	8
3190	CB	ASN	A	401	21.990	50.632	52.590	1.00	16.43	6
3191	CG	ASN	A	401	21.827	49.617	51.404	1.00	25.29	6
3192	OD1	ASN	A	401	22.805	49.256	50.797	1.00	33.03	8
3193	ND2	ASN	A	401	20.619	49.215	51.102	1.00	40.20	7
3194	N	ASN	A	402	21.705	53.948	53.373	1.00	13.76	7
3195	CA	ASN	A	402	21.449	54.851	54.506	1.00	13.71	6
3196	C	ASN	A	402	20.938	56.175	53.980	1.00	14.67	6
3197	O	ASN	A	402	21.601	56.923	55.311	1.00	13.53	8
3198	CB	ASN	A	402	22.753	55.036	55.311	1.00	13.24	6
3199	CG	ASN	A	402	22.397	55.744	56.604	1.00	14.53	6
3200	OD1	ASN	A	402	21.722	56.787	56.564	1.00	14.56	8
3201	ND2	ASN	A	402	22.839	55.186	57.762	1.00	11.48	7
3202	N	ALA	A	403	19.633	56.460	54.202	1.00	12.55	6
3203	CA	ALA	A	403	19.017	57.672	53.679	1.00	12.81	6
3204	C	ALA	A	403	19.622	58.961	54.144	1.00	14.88	6
3205	O	ALA	A	403	19.421	59.958	53.422	1.00	13.18	8
3206	CB	ALA	A	403	17.511	57.608	54.092	1.00	12.99	6
3207	N	ALA	A	404	20.407	58.945	55.229	1.00	11.82	7
3208	CA	ALA	A	404	21.107	60.187	55.611	1.00	11.98	6
3209	C	ALA	A	404	22.095	60.641	54.524	1.00	11.49	6
3210	O	ALA	A	404	22.259	61.857	54.383	1.00	12.58	8
3211	CB	ALA	A	404	21.874	60.008	56.930	1.00	15.02	6
3212	N	ILE	A	405	22.707	59.716	53.830	1.00	11.27	7
3213	CA	ILE	A	405	23.702	60.168	52.791	1.00	11.62	6
3214	C	ILE	A	405	22.936	60.722	51.581	1.00	13.21	6
3215	O	ILE	A	405	23.353	61.696	50.964	1.00	12.14	8
3216	CB	ILE	A	405	24.538	58.973	52.331	1.00	14.54	6
3217	CG1	ILE	A	405	25.425	58.253	53.392	1.00	15.40	6
3218	CG2	ILE	A	405	25.511	59.410	51.202	1.00	11.30	6
3219	CD1	ILE	A	405	26.170	59.266	54.247	1.00	17.36	6
3220	N	GLN	A	406	21.759	60.152	51.297	1.00	11.32	7
3221	CA	GLN	A	406	20.992	60.545	50.107	1.00	11.87	6
3222	C	GLN	A	406	20.335	61.902	50.300	1.00	14.28	6

Serial	Atom	Res	Chain	Seq	X	Y	Z	Occ	B	n
3223	O	GLN	A	406	20.265	62.707	49.360	1.00	11.82	8
3224	CB	GLN	A	406	19.875	59.494	49.869	1.00	12.00	6
3225	CG	GLN	A	406	20.511	58.136	49.549	1.00	12.17	6
3226	CD	GLN	A	406	19.521	56.983	49.392	1.00	22.22	6
3227	OE1	GLN	A	406	19.847	55.775	49.395	1.00	20.31	8
3228	NE2	GLN	A	406	18.272	57.356	49.223	1.00	24.04	7
3229	N	TYR	A	407	19.757	62.154	51.499	1.00	13.05	7
3230	CA	TYR	A	407	18.848	63.261	51.702	1.00	13.56	6
3231	C	TYR	A	407	19.168	64.182	52.885	1.00	10.35	6
3232	O	TYR	A	407	18.463	65.190	53.047	1.00	12.85	8
3233	CB	TYR	A	407	17.440	62.678	52.068	1.00	14.06	6
3234	CG	TYR	A	407	16.935	61.567	51.144	1.00	12.51	6
3235	CD1	TYR	A	407	16.929	61.748	49.759	1.00	11.68	6
3236	CD2	TYR	A	407	16.397	60.418	51.708	1.00	11.57	6
3237	CE1	TYR	A	407	16.441	60.742	48.918	1.00	12.71	6
3238	CE2	TYR	A	407	15.910	59.420	50.894	1.00	14.60	6
3239	CZ	TYR	A	407	15.933	59.595	49.515	1.00	15.50	6
3240	OH	TYR	A	407	15.433	58.598	48.707	1.00	15.89	8
3241	N	GLY	A	408	20.152	63.743	53.678	1.00	12.31	7
3242	CA	GLY	A	408	20.277	64.361	55.005	1.00	14.17	6
3243	C	GLY	A	408	20.830	65.790	55.030	1.00	13.76	6
3244	O	GLY	A	408	21.566	66.253	54.133	1.00	13.31	8
3245	N	THR	A	409	20.636	66.393	56.207	1.00	12.33	7
3246	CA	THR	A	409	21.322	67.370	56.558	1.00	14.50	6
3247	C	THR	A	409	22.822	67.645	56.707	1.00	13.72	6
3248	O	THR	A	409	23.222	66.229	56.777	1.00	12.71	8
3249	CB	THR	A	409	20.812	68.210	57.928	1.00	16.13	6
3250	OG1	THR	A	409	20.849	67.169	56.916	1.00	16.69	8
3251	CG2	THR	A	409	19.360	68.677	57.762	1.00	18.70	6
3252	N	THR	A	410	23.605	68.449	56.790	1.00	12.28	7
3253	CA	THR	A	410	25.062	68.261	57.141	1.00	12.15	6
3254	C	THR	A	410	25.332	69.229	58.321	1.00	12.71	6
3255	O	THR	A	410	25.041	70.423	58.225	1.00	13.99	8
3256	CB	THR	A	410	25.943	68.740	55.952	1.00	11.34	6
3257	OG1	THR	A	410	25.541	67.959	54.786	1.00	13.77	8
3258	CG2	THR	A	410	27.425	68.392	56.192	1.00	13.33	6
3259	N	THR	A	411	25.996	68.665	59.342	1.00	12.78	7
3260	CA	THR	A	411	26.301	69.430	60.568	1.00	12.27	6
3261	C	THR	A	411	27.765	69.192	60.919	1.00	13.51	6
3262	O	THR	A	411	28.168	68.039	61.036	1.00	14.10	8
3263	CB	THR	A	411	25.386	68.891	61.718	1.00	13.83	6
3264	OG1	THR	A	411	24.000	69.120	61.364	1.00	15.26	8
3265	CG2	THR	A	411	25.620	69.676	62.991	1.00	17.79	6
3266	N	GLN	A	412	28.501	70.302	61.110	1.00	12.74	7
3267	CA	GLN	A	412	29.899	70.153	61.592	1.00	13.61	6
3268	C	GLN	A	412	29.894	69.765	63.062	1.00	13.43	6
3269	O	GLN	A	412	29.218	70.451	63.824	1.00	15.64	8
3270	CB	GLN	A	412	30.556	71.523	61.335	1.00	14.70	6
3271	CG	GLN	A	412	31.999	71.615	61.876	1.00	22.99	6
3272	CD	GLN	A	412	31.918	72.168	63.334	1.00	22.07	6
3273	OE1	GLN	A	412	32.409	71.438	64.163	1.00	19.41	8
3274	NE2	GLN	A	412	31.358	73.310	63.633	1.00	25.10	7
3275	N	ARG	A	413	30.694	68.721	63.330	1.00	11.94	7
3276	CA	ARG	A	413	30.768	68.223	64.700	1.00	11.40	6
3277	C	ARG	A	413	32.184	68.430	65.276	1.00	12.19	6
3278	O	ARG	A	413	32.263	68.482	66.510	1.00	12.20	8
3279	CB	ARG	A	413	30.356	66.756	64.815	1.00	16.20	6
3280	CG	ARG	A	413	28.840	66.557	64.562	1.00	13.09	6
3281	CD	ARG	A	413	27.968	67.296	65.579	1.00	13.13	6
3282	NE	ARG	A	413	26.611	66.668	65.580	1.00	14.79	7
3283	CZ	ARG	A	413	25.684	66.909	66.515	1.00	19.43	6
3284	NH1	ARG	A	413	25.974	67.784	67.503	1.00	16.73	7
3285	NH2	ARG	A	413	24.529	66.233	66.470	1.00	18.22	7
3286	N	TRP	A	414	33.228	68.466	64.474	1.00	11.18	7
3287	CA	TRP	A	414	34.548	68.752	65.082	1.00	12.61	6
3288	C	TRP	A	414	35.444	69.243	63.947	1.00	13.80	6
3289	O	TRP	A	414	35.295	68.712	62.833	1.00	13.58	8
3290	CB	TRP	A	414	35.121	67.438	65.651	1.00	14.04	6
3291	CG	TRP	A	414	36.181	67.603	66.709	1.00	14.49	6
3292	CD1	TRP	A	414	35.921	67.473	68.075	1.00	15.45	6
3293	CD2	TRP	A	414	37.583	67.864	66.591	1.00	15.74	6
3294	NE1	TRP	A	414	37.112	67.622	68.773	1.00	16.17	7
3295	CE2	TRP	A	414	38.123	67.884	67.898	1.00	15.43	6
3296	CE3	TRP	A	414	38.421	68.101	65.507	1.00	14.89	6
3297	CZ2	TRP	A	414	39.489	68.090	68.138	1.00	20.21	6
3298	CZ3	TRP	A	414	39.805	68.279	65.725	1.00	15.42	6
3299	CH2	TRP	A	414	40.310	68.296	67.066	1.00	17.01	6
3300	N	ILE	A	415	36.215	70.337	64.186	1.00	12.40	7
3301	CA	ILE	A	415	37.012	70.828	63.038	1.00	11.33	6
3302	C	ILE	A	415	38.296	71.491	63.524	1.00	15.04	6
3303	O	ILE	A	415	38.300	72.134	64.600	1.00	15.39	8
3304	CB	ILE	A	415	36.153	71.827	62.253	1.00	15.29	6
3305	CG1	ILE	A	415	36.843	72.197	60.923	1.00	15.07	6
3306	CG2	ILE	A	415	35.782	73.061	63.054	1.00	20.69	6

3307	CD1	ILE	A	415	35.862	72.686	59.854	1.00	18.18	6
3308	N	ASN	A	416	39.349	71.266	62.797	1.00	13.47	7
3309	CA	ASN	A	416	40.517	72.152	62.912	1.00	14.63	6
3310	C	ASN	A	416	41.103	72.198	61.526	1.00	13.17	6
3311	O	ASN	A	416	40.399	71.879	60.536	1.00	12.76	8
3312	CB	ASN	A	416	41.480	71.607	63.994	1.00	14.61	6
3313	CG	ASN	A	416	42.250	70.394	63.669	1.00	16.07	6
3314	OD1	ASN	A	416	42.180	69.778	62.591	1.00	12.99	8
3315	ND2	ASN	A	416	43.129	69.947	64.597	1.00	16.12	7
3316	N	ASN	A	417	42.356	72.644	61.313	1.00	13.23	7
3317	CA	ASN	A	417	42.755	72.836	59.907	1.00	13.30	6
3318	C	ASN	A	417	42.767	71.526	59.103	1.00	13.34	6
3319	O	ASN	A	417	42.621	71.524	57.886	1.00	12.55	8
3320	CB	ASN	A	417	44.185	73.410	59.931	1.00	16.87	6
3321	CG	ASN	A	417	44.190	74.929	59.999	1.00	23.81	6
3322	OD1	ASN	A	417	43.170	75.585	60.178	1.00	21.65	8
3323	ND2	ASN	A	417	45.415	75.482	59.839	1.00	24.48	7
3324	N	ASP	A	418	43.109	70.439	59.874	1.00	12.62	7
3325	CA	ASP	A	418	43.295	69.150	59.209	1.00	10.76	6
3326	C	ASP	A	418	42.126	68.157	59.319	1.00	12.08	6
3327	O	ASP	A	418	42.093	67.177	58.533	1.00	11.98	8
3328	CB	ASP	A	418	44.520	68.478	59.846	1.00	13.38	6
3329	CG	ASP	A	418	45.842	69.080	59.483	1.00	16.78	6
3330	OD1	ASP	A	418	46.040	69.609	58.372	1.00	13.55	8
3331	OD2	ASP	A	418	46.759	69.017	60.342	1.00	13.98	8
3332	N	VAL	A	419	41.304	68.369	60.363	1.00	11.52	7
3333	CA	VAL	A	419	40.281	67.366	60.644	1.00	10.52	6
3334	C	VAL	A	419	38.910	67.958	60.419	1.00	12.12	6
3335	O	VAL	A	419	38.606	69.052	60.846	1.00	12.25	8
3336	CB	VAL	A	419	40.384	66.958	62.145	1.00	11.46	6
3337	CG1	VAL	A	419	39.231	65.959	62.481	1.00	11.88	6
3338	CG2	VAL	A	419	41.737	66.266	62.444	1.00	12.56	6
3339	N	TYR	A	420	38.034	67.172	59.764	1.00	11.92	7
3340	CA	TYR	A	420	36.628	67.548	59.636	1.00	10.72	6
3341	C	TYR	A	420	35.772	66.328	59.977	1.00	10.50	6
3342	O	TYR	A	420	35.908	65.254	59.383	1.00	13.20	8
3343	CB	TYR	A	420	36.410	67.866	58.138	1.00	9.44	6
3344	CG	TYR	A	420	35.063	68.583	57.856	1.00	10.31	6
3345	CD1	TYR	A	420	34.258	69.238	58.754	1.00	13.30	6
3346	CD2	TYR	A	420	34.669	68.523	56.512	1.00	13.25	6
3347	CE1	TYR	A	420	33.045	69.865	58.379	1.00	11.70	6
3348	CE2	TYR	A	420	33.476	69.131	56.115	1.00	14.40	6
3349	CZ	TYR	A	420	32.692	69.780	57.021	1.00	15.49	6
3350	OH	TYR	A	420	31.503	70.412	56.683	1.00	15.26	8
3351	N	ILE	A	421	35.000	66.508	61.075	1.00	11.61	7
3352	CA	ILE	A	421	34.037	65.462	61.458	1.00	9.13	6
3353	C	ILE	A	421	32.649	66.076	61.282	1.00	10.90	6
3354	O	ILE	A	421	32.379	67.148	61.788	1.00	13.13	8
3355	CB	ILE	A	421	34.238	64.997	62.941	1.00	11.03	6
3356	CG1	ILE	A	421	35.689	64.584	63.133	1.00	11.76	6
3357	CG2	ILE	A	421	33.251	63.881	63.198	1.00	14.22	6
3358	CD1	ILE	A	421	35.980	63.871	64.484	1.00	14.70	6
3359	N	TYR	A	422	31.889	65.498	60.276	1.00	10.51	7
3360	CA	TYR	A	422	30.571	66.072	60.004	1.00	11.77	6
3361	C	TYR	A	422	29.510	64.963	60.053	1.00	11.33	6
3362	O	TYR	A	422	29.877	63.774	59.986	1.00	11.55	8
3363	CB	TYR	A	422	30.570	66.791	58.623	1.00	11.55	6
3364	CG	TYR	A	422	31.000	65.881	57.458	1.00	12.40	6
3365	CD1	TYR	A	422	30.094	65.103	56.724	1.00	11.81	6
3366	CD2	TYR	A	422	32.354	65.851	57.140	1.00	11.72	6
3367	CE1	TYR	A	422	30.559	64.305	55.668	1.00	12.09	6
3368	CE2	TYR	A	422	32.853	65.051	56.094	1.00	11.73	6
3369	CZ	TYR	A	422	31.935	64.298	55.377	1.00	13.04	6
3370	OH	TYR	A	422	32.365	63.520	54.338	1.00	11.43	8
3371	N	GLU	A	423	28.257	65.380	60.139	1.00	11.98	7
3372	CA	GLU	A	423	27.202	64.378	60.362	1.00	11.04	6
3373	C	GLU	A	423	26.064	64.598	59.383	1.00	11.74	6
3374	O	GLU	A	423	25.644	65.748	59.205	1.00	13.56	8
3375	CB	GLU	A	423	26.633	64.609	61.806	1.00	14.91	6
3376	CG	GLU	A	423	25.731	63.465	62.258	1.00	12.27	6
3377	CD	GLU	A	423	25.459	63.688	63.796	1.00	13.28	6
3378	OE1	GLU	A	423	24.947	64.750	64.099	1.00	18.28	8
3379	OE2	GLU	A	423	25.800	62.682	64.408	1.00	17.49	8
3380	N	ARG	A	424	25.648	63.541	58.699	1.00	12.31	7
3381	CA	ARG	A	424	24.457	63.621	57.847	1.00	11.64	6
3382	C	ARG	A	424	23.268	63.035	58.643	1.00	13.19	6
3383	O	ARG	A	424	23.515	62.077	59.367	1.00	13.04	8
3384	CB	ARG	A	424	24.665	62.689	56.620	1.00	10.72	6
3385	CG	ARG	A	424	25.961	63.049	55.805	1.00	10.81	6
3386	CD	ARG	A	424	25.862	64.465	55.212	1.00	11.09	6
3387	NE	ARG	A	424	24.666	64.721	54.412	1.00	10.91	7
3388	CZ	ARG	A	424	24.420	64.168	53.207	1.00	11.39	6
3389	NH1	ARG	A	424	25.240	63.308	52.620	1.00	10.62	7
3390	NH2	ARG	A	424	23.239	64.512	52.596	1.00	10.78	7

Atom#	Atom	Res	Chain	ResNum	X	Y	Z	Occ	B	
3391	N	LYS	A	425	22.068	63.623	58.542	1.00	11.90	7
3392	CA	LYS	A	425	20.937	63.080	59.290	1.00	11.75	6
3393	C	LYS	A	425	19.685	63.233	58.443	1.00	15.39	6
3394	O	LYS	A	425	19.444	64.340	58.036	1.00	14.23	8
3395	CB	LYS	A	425	20.801	63.854	60.626	1.00	13.68	6
3396	CG	LYS	A	425	19.617	63.271	61.494	1.00	14.44	6
3397	CD	LYS	A	425	19.721	63.992	62.863	1.00	19.53	6
3398	CE	LYS	A	425	18.739	63.342	63.856	1.00	20.97	6
3399	NZ	LYS	A	425	17.322	63.663	63.509	1.00	23.72	7
3400	N	PHE	A	426	18.908	62.144	58.357	1.00	12.33	7
3401	CA	PHE	A	426	17.583	62.238	57.709	1.00	12.03	6
3402	C	PHE	A	426	16.651	61.481	58.706	1.00	11.05	6
3403	O	PHE	A	426	16.657	60.252	58.693	1.00	13.03	8
3404	CB	PHE	A	426	17.608	61.450	56.397	1.00	13.23	6
3405	CG	PHE	A	426	16.260	61.449	55.699	1.00	12.58	6
3406	CD1	PHE	A	426	15.838	62.659	55.147	1.00	14.04	6
3407	CD2	PHE	A	426	15.515	60.300	55.571	1.00	14.55	6
3408	CE1	PHE	A	426	14.594	62.715	54.490	1.00	13.58	6
3409	CE2	PHE	A	426	14.264	60.345	54.906	1.00	15.67	6
3410	CZ	PHE	A	426	13.839	61.552	54.398	1.00	15.62	6
3411	N	PHE	A	427	15.916	62.357	59.408	1.00	14.34	7
3412	CA	PHE	A	427	15.024	61.773	60.490	1.00	16.20	6
3413	C	PHE	A	427	15.781	60.836	61.400	1.00	17.32	6
3414	O	PHE	A	427	16.737	61.354	62.070	1.00	19.10	8
3415	CB	PHE	A	427	13.772	61.132	59.840	1.00	18.52	6
3416	CG	PHE	A	427	12.888	62.175	59.193	1.00	17.73	6
3417	CD1	PHE	A	427	11.972	62.918	59.905	1.00	21.55	6
3418	CD2	PHE	A	427	13.018	62.396	57.830	1.00	14.51	6
3419	CE1	PHE	A	427	11.188	63.858	59.276	1.00	18.85	6
3420	CE2	PHE	A	427	12.246	63.336	57.185	1.00	16.68	6
3421	CZ	PHE	A	427	11.311	64.087	57.906	1.00	19.85	6
3422	N	ASN	A	428	15.546	59.522	61.442	1.00	15.71	7
3423	CA	ASN	A	428	16.284	58.735	62.433	1.00	17.74	6
3424	C	ASN	A	428	17.573	58.173	61.837	1.00	18.72	6
3425	O	ASN	A	428	18.307	57.587	62.639	1.00	17.57	8
3426	CB	ASN	A	428	15.484	57.591	63.008	1.00	22.35	6
3427	CG	ASN	A	428	14.267	58.214	63.733	1.00	36.36	6
3428	OD1	ASN	A	428	14.452	59.092	64.570	1.00	38.45	8
3429	ND2	ASN	A	428	13.103	63.325	63.325	1.00	44.64	7
3430	N	ASP	A	429	17.867	58.370	60.547	1.00	15.06	7
3431	CA	ASP	A	429	19.091	57.810	59.986	1.00	14.19	6
3432	C	ASP	A	429	20.222	58.816	60.103	1.00	11.55	6
3433	O	ASP	A	429	20.026	60.024	59.967	1.00	13.16	8
3434	CB	ASP	A	429	18.927	57.501	58.471	1.00	12.81	6
3435	CG	ASP	A	429	17.766	56.561	58.255	1.00	28.25	6
3436	OD1	ASP	A	429	17.727	55.530	58.940	1.00	20.84	8
3437	OD2	ASP	A	429	16.804	56.870	57.490	1.00	27.47	8
3438	N	VAL	A	430	21.384	58.297	60.600	1.00	12.85	7
3439	CA	VAL	A	430	22.502	59.192	60.876	1.00	11.51	6
3440	C	VAL	A	430	23.816	58.621	60.371	1.00	13.31	6
3441	O	VAL	A	430	23.993	57.405	60.522	1.00	12.49	8
3442	CB	VAL	A	430	22.632	59.315	62.427	1.00	14.21	6
3443	CG1	VAL	A	430	23.828	60.233	62.734	1.00	14.56	6
3444	CG2	VAL	A	430	21.373	59.955	63.033	1.00	14.79	6
3445	N	VAL	A	431	24.721	59.432	59.826	1.00	12.54	7
3446	CA	VAL	A	431	26.043	58.933	59.434	1.00	11.58	6
3447	C	VAL	A	431	27.051	59.978	59.912	1.00	11.43	6
3448	O	VAL	A	431	26.847	61.129	59.603	1.00	12.12	8
3449	CB	VAL	A	431	26.250	58.685	57.905	1.00	12.54	6
3450	CG1	VAL	A	431	27.698	58.199	57.615	1.00	10.81	6
3451	CG2	VAL	A	431	25.254	57.657	57.400	1.00	13.14	6
3452	N	LEU	A	432	27.976	59.557	60.800	1.00	10.48	7
3453	CA	LEU	A	432	29.015	60.512	61.285	1.00	11.27	6
3454	C	LEU	A	432	30.317	60.140	60.583	1.00	11.12	6
3455	O	LEU	A	432	30.708	58.961	60.577	1.00	13.28	8
3456	CB	LEU	A	432	29.087	60.224	62.831	1.00	11.86	6
3457	CG	LEU	A	432	30.033	61.205	63.562	1.00	11.74	6
3458	CD1	LEU	A	432	29.574	62.643	63.510	1.00	11.18	6
3459	CD2	LEU	A	432	30.119	60.744	65.044	1.00	14.23	6
3460	N	VAL	A	433	30.987	61.141	59.978	1.00	11.73	7
3461	CA	VAL	A	433	32.162	60.867	59.131	1.00	9.07	6
3462	C	VAL	A	433	33.328	61.667	59.702	1.00	9.49	6
3463	O	VAL	A	433	33.158	62.864	59.844	1.00	10.58	8
3464	CB	VAL	A	433	31.899	61.334	57.674	1.00	10.62	6
3465	CG1	VAL	A	433	33.173	61.088	56.812	1.00	11.05	6
3466	CG2	VAL	A	433	30.693	60.564	57.072	1.00	11.15	6
3467	N	ALA	A	434	34.455	61.000	60.017	1.00	10.65	7
3468	CA	ALA	A	434	35.643	61.757	60.409	1.00	10.07	6
3469	C	ALA	A	434	36.742	61.603	59.363	1.00	11.73	6
3470	O	ALA	A	434	37.030	60.489	58.924	1.00	11.96	8
3471	CB	ALA	A	434	36.199	61.171	61.742	1.00	10.24	6
3472	N	ILE	A	435	37.345	62.744	62.992	1.00	10.23	7
3473	CA	ILE	A	435	38.438	62.741	58.008	1.00	8.82	6
3474	C	ILE	A	435	39.571	63.587	58.558	1.00	10.58	6

Serial	Atom	Res	Chain	Seq	X	Y	Z	Occ	B	Elem
3475	O	ILE	A	435	39.366	64.734	58.922	1.00	10.88	8
3476	CB	ILE	A	435	37.961	63.408	56.674	1.00	10.84	6
3477	CG1	ILE	A	435	36.749	62.608	56.120	1.00	11.49	6
3478	CG2	ILE	A	435	39.146	63.352	55.689	1.00	10.37	6
3479	CD1	ILE	A	435	36.230	63.228	54.758	1.00	13.59	6
3480	N	ASN	A	436	40.741	62.933	58.596	1.00	10.83	7
3481	CA	ASN	A	436	41.963	63.646	58.943	1.00	11.64	6
3482	C	ASN	A	436	42.852	63.653	57.715	1.00	9.83	6
3483	O	ASN	A	436	43.367	62.621	57.309	1.00	12.85	8
3484	CB	ASN	A	436	42.711	62.821	60.036	1.00	12.59	6
3485	CG	ASN	A	436	44.030	63.480	60.383	1.00	12.87	6
3486	OD1	ASN	A	436	44.422	64.602	60.026	1.00	13.07	8
3487	ND2	ASN	A	436	44.789	62.721	61.224	1.00	13.56	7
3488	N	ARG	A	437	43.055	64.891	57.181	1.00	10.03	7
3489	CA	ARG	A	437	43.878	64.887	55.958	1.00	9.39	6
3490	C	ARG	A	437	45.362	64.725	56.267	1.00	10.71	6
3491	O	ARG	A	437	46.117	64.552	55.310	1.00	12.32	8
3492	CB	ARG	A	437	43.673	66.201	55.161	1.00	12.80	6
3493	CG	ARG	A	437	44.296	67.415	55.869	1.00	13.58	6
3494	CD	ARG	A	437	44.031	68.719	55.043	1.00	11.13	6
3495	NE	ARG	A	437	44.772	69.847	55.738	1.00	10.90	7
3496	CZ	ARG	A	437	44.957	71.046	55.195	1.00	14.88	6
3497	NH1	ARG	A	437	44.521	71.299	53.948	1.00	11.67	7
3498	NH2	ARG	A	437	45.406	72.058	55.947	1.00	13.84	7
3499	N	ASN	A	438	45.779	65.035	57.513	1.00	11.11	7
3500	CA	ASN	A	438	47.255	65.048	57.734	1.00	12.20	6
3501	C	ASN	A	438	47.780	63.623	57.869	1.00	13.73	6
3502	O	ASN	A	438	47.440	62.900	58.830	1.00	12.56	8
3503	CB	ASN	A	438	47.474	65.787	59.071	1.00	13.49	6
3504	CG	ASN	A	438	48.921	66.167	59.255	1.00	13.92	6
3505	OD1	ASN	A	438	49.263	65.335	58.985	1.00	15.45	8
3506	ND2	ASN	A	438	49.263	67.376	59.700	1.00	13.51	7
3507	N	THR	A	439	48.650	63.224	56.949	1.00	11.49	7
3508	CA	THR	A	439	49.087	61.822	56.883	1.00	12.77	6
3509	C	THR	A	439	50.271	61.585	57.846	1.00	14.69	6
3510	O	THR	A	439	50.750	60.450	57.917	1.00	16.23	8
3511	CB	THR	A	439	49.518	61.431	55.456	1.00	16.48	6
3512	OG1	THR	A	439	50.542	62.330	54.990	1.00	17.73	8
3513	CG2	THR	A	439	48.284	61.611	54.554	1.00	15.47	6
3514	N	GLN	A	440	50.666	62.669	58.501	1.00	14.00	7
3515	CA	GLN	A	440	51.778	62.467	59.481	1.00	16.61	6
3516	C	GLN	A	440	51.339	62.729	60.910	1.00	19.80	6
3517	O	GLN	A	440	52.198	62.543	61.812	1.00	18.55	8
3518	CB	GLN	A	440	52.905	63.422	59.081	1.00	23.60	6
3519	CG	GLN	A	440	53.543	62.971	57.761	1.00	37.17	6
3520	CD	GLN	A	440	54.248	61.634	57.810	1.00	48.48	6
3521	OE1	GLN	A	440	55.292	61.462	58.444	1.00	57.64	8
3522	NE2	GLN	A	440	53.729	60.605	57.127	1.00	54.59	7
3523	N	SER	A	441	50.115	63.161	61.192	1.00	13.09	7
3524	CA	SER	A	441	49.738	63.516	62.559	1.00	13.14	6
3525	C	SER	A	441	48.481	62.739	62.958	1.00	18.60	6
3526	O	SER	A	441	47.524	62.718	62.168	1.00	15.27	8
3527	CB	SER	A	441	49.451	65.025	62.702	1.00	16.00	6
3528	OG	SER	A	441	50.658	65.770	62.516	1.00	17.34	8
3529	N	SER	A	442	48.417	62.251	64.198	1.00	13.46	7
3530	CA	SER	A	442	47.177	61.703	64.730	1.00	13.64	6
3531	C	SER	A	442	46.581	62.835	65.599	1.00	14.33	6
3532	O	SER	A	442	47.366	63.761	65.974	1.00	15.22	8
3533	CB	SER	A	442	47.452	60.598	65.791	1.00	15.48	6
3534	OG	SER	A	442	48.033	59.420	65.203	1.00	19.49	8
3535	N	TYR	A	443	45.302	62.845	65.817	1.00	13.24	7
3536	CA	TYR	A	443	44.693	63.886	66.694	1.00	12.26	6
3537	C	TYR	A	443	43.877	63.156	67.749	1.00	14.89	6
3538	O	TYR	A	443	43.032	62.293	67.451	1.00	14.44	8
3539	CB	TYR	A	443	43.772	64.824	65.860	1.00	13.59	6
3540	CG	TYR	A	443	44.592	65.807	65.009	1.00	10.92	6
3541	CD1	TYR	A	443	45.124	66.940	65.629	1.00	12.95	6
3542	CD2	TYR	A	443	44.836	65.549	63.667	1.00	12.89	6
3543	CE1	TYR	A	443	45.908	67.843	64.882	1.00	11.89	6
3544	CE2	TYR	A	443	45.595	66.474	62.914	1.00	11.54	6
3545	CZ	TYR	A	443	46.107	67.595	63.551	1.00	12.16	6
3546	OH	TYR	A	443	46.871	68.508	62.811	1.00	15.42	8
3547	N	SER	A	444	43.976	63.663	69.008	1.00	14.60	7
3548	CA	SER	A	444	43.034	63.232	70.040	1.00	14.67	6
3549	C	SER	A	444	41.698	64.002	69.912	1.00	13.95	6
3550	O	SER	A	444	41.739	65.229	69.879	1.00	16.35	8
3551	CB	SER	A	444	43.620	63.684	71.418	1.00	17.27	6
3552	OG	SER	A	444	44.701	62.764	71.733	1.00	18.88	8
3553	N	ILE	A	445	40.628	63.265	69.758	1.00	15.00	7
3554	CA	ILE	A	445	39.303	63.892	69.559	1.00	13.10	6
3555	C	ILE	A	445	38.498	63.757	70.863	1.00	14.11	6
3556	O	ILE	A	445	37.935	62.703	71.131	1.00	17.51	8
3557	CB	ILE	A	445	38.537	63.188	68.403	1.00	14.59	6
3558	CG1	ILE	A	445	39.386	63.120	67.136	1.00	15.39	6

Serial	Atom	Res	Chain	ResNum	X	Y	Z	Occ	B	El
3559	CG2	ILE	A	445	37.159	63.906	68.223	1.00	16.44	6
3560	CD1	ILE	A	445	39.772	64.500	66.533	1.00	15.53	6
3561	N	SER	A	446	38.309	64.925	71.501	1.00	15.73	7
3562	CA	SER	A	446	37.450	65.018	72.680	1.00	16.64	6
3563	C	SER	A	446	36.394	66.091	72.415	1.00	18.39	6
3564	O	SER	A	446	36.592	67.021	71.653	1.00	18.43	8
3565	CB	SER	A	446	38.248	65.398	73.972	1.00	18.77	6
3566	OG	SER	A	446	38.784	66.689	73.750	1.00	24.38	8
3567	N	GLY	A	447	35.263	65.958	73.091	1.00	19.05	7
3568	CA	GLY	A	447	34.169	66.916	73.050	1.00	17.05	6
3569	C	GLY	A	447	33.253	66.660	71.829	1.00	20.26	6
3570	O	GLY	A	447	32.491	67.554	71.467	1.00	19.38	8
3571	N	LEU	A	448	33.487	65.502	71.171	1.00	13.94	7
3572	CA	LEU	A	448	32.625	65.262	69.967	1.00	15.40	6
3573	C	LEU	A	448	31.245	64.853	70.386	1.00	16.78	6
3574	O	LEU	A	448	31.018	63.906	71.155	1.00	17.47	8
3575	CB	LEU	A	448	33.312	64.155	69.144	1.00	14.08	6
3576	CG	LEU	A	448	32.578	63.628	67.903	1.00	15.34	6
3577	CD1	LEU	A	448	32.403	64.740	66.845	1.00	16.18	6
3578	CD2	LEU	A	448	33.283	62.449	67.256	1.00	13.42	6
3579	N	GLN	A	449	30.221	65.509	69.814	1.00	13.16	7
3580	CA	GLN	A	449	28.812	65.224	70.018	1.00	14.91	6
3581	C	GLN	A	449	28.209	64.630	68.734	1.00	16.76	6
3582	O	GLN	A	449	28.754	64.827	67.645	1.00	13.94	8
3583	CB	AGLN	A	449	28.054	66.486	70.470	0.50	18.23	6
3584	CG	AGLN	A	449	28.884	67.206	71.540	0.50	24.01	6
3585	CD	AGLN	A	449	28.211	68.188	72.449	0.50	23.44	6
3586	OE1	AGLN	A	449	28.812	68.606	73.455	0.50	22.43	8
3587	NE2	AGLN	A	449	26.984	68.573	72.134	0.50	32.22	7
3583	CB	BGLN	A	449	27.974	66.473	70.339	0.50	13.26	6
3584	CG	BGLN	A	449	28.536	67.122	71.620	0.50	17.72	6
3585	CD	BGLN	A	449	28.037	66.396	72.844	0.50	16.00	6
3586	OE1	BGLN	A	449	28.776	65.702	73.511	0.50	22.43	8
3587	NE2	BGLN	A	449	26.759	66.550	73.145	0.50	23.34	7
3588	N	THR	A	450	27.085	63.965	68.897	1.00	15.61	7
3589	CA	THR	A	450	26.481	63.264	67.761	1.00	13.97	6
3590	C	THR	A	450	24.998	63.001	67.994	1.00	16.43	6
3591	O	THR	A	450	24.528	62.925	69.160	1.00	16.61	8
3592	CB	THR	A	450	27.216	61.912	67.566	1.00	15.00	6
3593	OG1	THR	A	450	26.686	61.176	66.471	1.00	15.07	8
3594	CG2	THR	A	450	27.007	61.016	68.814	1.00	18.27	6
3595	N	ALA	A	451	24.286	62.806	66.886	1.00	15.10	7
3596	CA	ALA	A	451	22.898	62.330	66.941	1.00	16.96	6
3597	C	ALA	A	451	22.799	60.830	66.879	1.00	15.04	6
3598	O	ALA	A	451	21.681	60.282	66.900	1.00	15.44	8
3599	CB	ALA	A	451	22.212	62.946	65.681	1.00	16.26	6
3600	N	LEU	A	452	23.929	60.083	66.788	1.00	14.57	7
3601	CA	LEU	A	452	23.803	58.623	66.814	1.00	12.45	6
3602	C	LEU	A	452	23.170	58.200	68.158	1.00	17.32	6
3603	O	LEU	A	452	23.505	58.727	69.218	1.00	17.00	8
3604	CB	LEU	A	452	25.201	57.989	66.779	1.00	14.27	6
3605	CG	LEU	A	452	25.854	58.082	65.380	1.00	14.43	6
3606	CD1	LEU	A	452	27.344	57.716	65.610	1.00	13.81	6
3607	CD2	LEU	A	452	25.263	57.102	64.375	1.00	14.40	6
3608	N	PRO	A	453	22.383	57.163	68.080	1.00	16.12	7
3609	CA	PRO	A	453	21.808	56.539	69.305	1.00	18.87	6
3610	C	PRO	A	453	22.936	55.882	70.090	1.00	19.68	6
3611	O	PRO	A	453	24.039	55.532	69.686	1.00	18.19	8
3612	CB	PRO	A	453	20.796	55.523	68.880	1.00	20.41	6
3613	CG	PRO	A	453	20.615	55.743	67.392	1.00	21.35	6
3614	CD	PRO	A	453	21.806	56.582	66.872	1.00	16.78	6
3615	N	ASN	A	454	22.619	55.684	71.406	1.00	14.58	7
3616	CA	ASN	A	454	23.630	55.025	72.244	1.00	14.89	6
3617	C	ASN	A	454	24.085	53.704	71.712	1.00	18.47	6
3618	O	ASN	A	454	23.320	52.881	71.168	1.00	17.84	8
3619	CB	ASN	A	454	22.851	54.705	73.573	1.00	17.80	6
3620	CG	ASN	A	454	22.656	55.933	74.421	1.00	24.08	6
3621	OD1	ASN	A	454	23.071	57.055	74.201	1.00	19.24	8
3622	ND2	ASN	A	454	21.941	55.757	75.554	1.00	24.42	7
3623	N	GLY	A	455	25.378	53.451	71.950	1.00	17.75	7
3624	CA	GLY	A	455	25.919	52.156	71.553	1.00	20.13	6
3625	C	GLY	A	455	27.422	52.289	71.172	1.00	17.39	6
3626	O	GLY	A	455	27.899	53.393	71.080	1.00	19.07	8
3627	N	SER	A	456	27.916	51.112	70.814	1.00	18.37	7
3628	CA	SER	A	456	29.286	51.073	70.261	1.00	17.34	6
3629	C	SER	A	456	29.173	50.927	68.725	1.00	17.80	6
3630	O	SER	A	456	28.322	50.199	68.220	1.00	20.53	8
3631	CB	SER	A	456	29.916	49.747	70.778	1.00	22.39	6
3632	OG	SER	A	456	30.178	49.998	72.161	1.00	30.42	8
3633	N	TYR	A	457	30.024	51.683	68.013	1.00	15.08	7
3634	CA	TYR	A	457	29.941	51.627	66.559	1.00	13.18	6
3635	C	TYR	A	457	31.301	51.229	65.961	1.00	14.20	6
3636	O	TYR	A	457	32.257	51.915	66.259	1.00	16.66	8
3637	CB	TYR	A	457	29.564	53.017	65.941	1.00	17.47	6

Atom	Name	Residue	Chain	Seq	X	Y	Z	Occ	B	
3638	CG	TYR	A	457	28.122	53.387	66.241	1.00	15.26	6
3639	CD1	TYR	A	457	27.799	53.933	67.497	1.00	15.37	6
3640	CD2	TYR	A	457	27.077	53.174	65.325	1.00	16.16	6
3641	CE1	TYR	A	457	26.521	54.297	67.873	1.00	15.82	6
3642	CE2	TYR	A	457	25.768	53.516	65.671	1.00	15.99	6
3643	CZ	TYR	A	457	25.523	54.070	66.928	1.00	16.37	6
3644	OH	TYR	A	457	24.210	54.401	67.218	1.00	16.11	8
3645	N	ALA	A	458	31.329	50.211	65.128	1.00	14.59	7
3646	CA	ALA	A	458	32.601	49.943	64.445	1.00	17.03	6
3647	C	ALA	A	458	32.686	50.915	63.247	1.00	14.66	6
3648	O	ALA	A	458	31.654	51.326	62.731	1.00	14.84	8
3649	CB	ALA	A	458	32.638	48.538	63.857	1.00	17.50	6
3650	N	ASP	A	459	33.942	51.128	62.831	1.00	13.42	7
3651	CA	ASP	A	459	34.082	51.858	61.521	1.00	11.91	6
3652	C	ASP	A	459	33.472	51.048	60.410	1.00	12.75	6
3653	O	ASP	A	459	33.679	49.834	60.222	1.00	13.42	8
3654	CB	ASP	A	459	35.567	52.007	61.258	1.00	12.66	6
3655	CG	ASP	A	459	35.984	52.522	59.876	1.00	13.56	6
3656	OD1	ASP	A	459	35.143	53.232	59.328	1.00	12.36	8
3657	OD2	ASP	A	459	37.109	52.170	59.441	1.00	14.11	8
3658	N	TYR	A	460	32.581	51.730	59.662	1.00	11.96	7
3659	CA	TYR	A	460	31.927	51.022	58.513	1.00	13.29	6
3660	C	TYR	A	460	32.905	50.541	57.467	1.00	16.04	6
3661	O	TYR	A	460	32.617	49.627	56.683	1.00	15.32	8
3662	CB	TYR	A	460	30.909	52.021	57.931	1.00	13.60	6
3663	CG	TYR	A	460	29.970	51.418	56.899	1.00	13.06	6
3664	CD1	TYR	A	460	28.809	50.815	57.400	1.00	11.63	6
3665	CD2	TYR	A	460	30.215	51.365	55.532	1.00	16.49	6
3666	CE1	TYR	A	460	27.838	50.274	56.514	1.00	15.86	6
3667	CE2	TYR	A	460	29.278	50.783	54.662	1.00	15.66	6
3668	CZ	TYR	A	460	28.098	50.257	55.165	1.00	17.45	6
3669	OH	TYR	A	460	27.209	49.696	54.262	1.00	16.44	8
3670	N	LEU	A	461	34.057	51.242	57.374	1.00	13.04	7
3671	CA	LEU	A	461	35.137	50.814	56.424	1.00	13.89	6
3672	C	LEU	A	461	36.026	49.742	57.023	1.00	14.71	6
3673	O	LEU	A	461	36.992	49.331	56.369	1.00	13.39	8
3674	CB	LEU	A	461	35.968	52.115	56.171	1.00	13.08	6
3675	CG	LEU	A	461	35.299	53.075	55.149	1.00	13.44	6
3676	CD1	LEU	A	461	35.968	54.453	55.284	1.00	14.44	6
3677	CD2	LEU	A	461	35.485	52.526	53.743	1.00	16.37	7
3678	N	SER	A	462	35.746	49.218	58.222	1.00	15.11	7
3679	CA	SER	A	462	36.521	48.122	58.810	1.00	16.17	6
3680	C	SER	A	462	38.011	48.422	58.907	1.00	17.10	6
3681	O	SER	A	462	38.878	47.546	58.763	1.00	16.40	8
3682	CB	SER	A	462	36.316	46.813	57.987	1.00	19.17	6
3683	OG	SER	A	462	34.914	46.479	57.974	1.00	20.76	8
3684	N	GLY	A	463	38.369	49.689	57.115	1.00	13.98	7
3685	CA	GLY	A	463	39.781	50.089	59.115	1.00	14.97	6
3686	C	GLY	A	463	40.533	50.140	59.271	1.00	15.42	6
3687	O	GLY	A	463	41.745	50.387	57.949	1.00	15.84	8
3688	N	GLY	A	463	39.816	50.043	57.984	1.00	12.86	7
3689	CA	LEU	A	464	40.559	50.113	56.808	1.00	10.93	6
3690	C	LEU	A	464	41.370	51.420	55.525	1.00	12.62	6
3691	O	LEU	A	464	42.484	51.408	55.427	1.00	14.36	8
3692	CB	LEU	A	464	39.487	50.148	54.854	1.00	13.24	6
3693	CG	LEU	A	464	40.083	50.223	54.402	1.00	14.42	6
3694	CD1	LEU	A	464	40.800	48.892	52.969	1.00	15.93	6
3695	CD2	LEU	A	464	38.971	50.469	52.620	1.00	14.00	8
3696	N	LEU	A	465	40.797	52.526	51.967	1.00	11.79	7
3697	CA	LEU	A	465	41.473	53.824	55.872	1.00	12.51	6
3698	C	LEU	A	465	41.953	54.319	55.761	1.00	13.32	6
3699	O	LEU	A	465	41.873	55.518	57.114	1.00	16.12	8
3700	CB	LEU	A	465	40.510	54.894	57.401	1.00	12.90	6
3701	CG	LEU	A	465	40.090	54.441	55.130	1.00	12.88	6
3702	CD1	LEU	A	465	39.162	55.509	53.714	1.00	14.96	6
3703	CD2	LEU	A	465	41.263	54.191	53.111	1.00	14.83	8
3704	N	GLY	A	466	42.237	53.348	54.191	1.00	14.06	7
3705	CA	GLY	A	466	42.789	53.800	58.043	1.00	14.74	6
3706	C	GLY	A	466	41.735	54.095	59.336	1.00	14.06	6
3707	O	GLY	A	466	42.061	54.643	60.400	1.00	15.13	8
3708	N	GLY	A	467	40.451	53.817	61.479	1.00	13.86	7
3709	CA	GLY	A	467	39.357	54.137	60.125	1.00	11.80	6
3710	C	GLY	A	467	39.321	53.244	61.045	1.00	14.35	6
3711	O	GLY	A	467	40.083	52.261	62.311	1.00	16.59	8
3712	N	ASN	A	468	38.509	53.696	62.394	1.00	15.29	7
3713	CA	ASN	A	468	38.483	53.050	63.236	1.00	14.61	6
3714	C	ASN	A	468	37.007	52.960	64.569	1.00	15.14	6
3715	O	ASN	A	468	36.145	53.679	64.963	1.00	15.07	8
3716	CB	ASN	A	468	39.253	54.012	64.418	1.00	16.45	6
3717	CG	ASN	A	468	38.730	55.429	65.515	1.00	17.15	6
3718	OD1	ASN	A	468	39.013	56.375	65.538	1.00	19.92	8
3719	ND2	ASN	A	468	37.812	55.710	66.490	1.00	16.39	7
3720	N	GLY	A	469	36.787	52.285	66.076	1.00	16.49	7
3721	CA	GLY	A	469	35.415	52.189	66.624	1.00	16.85	6

#	Atom	Res	Ch	Seq	X	Y	Z	Occ	B	
3722	C	GLY	A	469	35.261	53.309	67.640	1.00	14.55	6
3723	O	GLY	A	469	36.191	53.919	68.176	1.00	17.51	8
3724	N	ILE	A	470	33.976	53.624	67.954	1.00	14.83	7
3725	CA	ILE	A	470	33.623	54.650	68.921	1.00	14.09	6
3726	C	ILE	A	470	32.518	54.102	69.860	1.00	13.92	6
3727	O	ILE	A	470	31.867	53.108	69.553	1.00	19.57	8
3728	CB	ILE	A	470	33.155	56.004	68.368	1.00	16.37	6
3729	CG1	ILE	A	470	31.948	55.753	67.429	1.00	17.51	6
3730	CG2	ILE	A	470	34.319	56.652	67.581	1.00	16.76	6
3731	CD1	ILE	A	470	31.315	57.104	67.032	1.00	15.78	6
3732	N	SER	A	471	32.408	54.804	70.980	1.00	15.71	7
3733	CA	SER	A	471	31.268	54.477	71.903	1.00	14.51	6
3734	C	SER	A	471	30.497	55.755	72.165	1.00	14.83	6
3735	O	SER	A	471	31.107	56.814	72.402	1.00	17.55	8
3736	CB	SER	A	471	32.024	54.018	73.220	1.00	22.00	8
3737	OG	SER	A	471	30.967	53.933	74.176	1.00	27.40	8
3738	N	VAL	A	472	29.172	55.712	72.054	1.00	14.69	7
3739	CA	VAL	A	472	28.298	56.834	72.186	1.00	14.15	6
3740	C	VAL	A	472	27.380	56.696	73.425	1.00	15.07	6
3741	O	VAL	A	472	26.764	55.638	73.563	1.00	19.05	8
3742	CB	VAL	A	472	27.433	56.903	70.900	1.00	15.53	6
3743	CG1	VAL	A	472	26.446	58.057	70.970	1.00	14.81	6
3744	CG2	VAL	A	472	28.376	57.153	69.691	1.00	18.40	8
3745	N	SER	A	473	27.351	57.767	74.184	1.00	15.80	7
3746	CA	SER	A	473	26.386	57.736	75.312	1.00	16.44	6
3747	C	SER	A	473	25.711	59.073	75.420	1.00	16.40	6
3748	O	SER	A	473	26.435	60.084	75.548	1.00	19.98	8
3749	CB	SER	A	473	27.214	57.446	76.594	1.00	21.00	6
3750	OG	SER	A	473	26.284	58.057	77.686	1.00	31.70	8
3751	N	ASN	A	474	24.396	59.107	75.308	1.00	18.40	7
3752	CA	ASN	A	474	23.632	60.348	75.444	1.00	19.45	6
3753	C	ASN	A	474	24.194	61.523	74.646	1.00	19.99	6
3754	O	ASN	A	474	24.363	62.658	75.103	1.00	15.86	8
3755	CB	ASN	A	474	23.685	60.707	76.968	1.00	22.35	6
3756	CG	ASN	A	474	22.991	59.572	77.722	1.00	31.24	6
3757	OD1	ASN	A	474	22.086	58.887	77.191	1.00	33.57	8
3758	ND2	ASN	A	474	23.579	59.303	78.890	1.00	32.67	7
3759	N	GLY	A	475	24.452	61.223	73.341	1.00	17.78	7
3760	CA	GLY	A	475	24.866	62.311	72.438	1.00	20.56	6
3761	C	GLY	A	475	26.369	62.573	72.430	1.00	18.34	6
3762	O	GLY	A	475	26.807	63.464	71.690	1.00	20.59	8
3763	N	SER	A	476	27.146	61.894	73.305	1.00	18.39	7

#	Atom	Res	Ch	Seq	X	Y	Z	Occ	B	
3764	CA	SER	A	476	28.563	62.222	73.451	1.00	14.26	6
3765	C	SER	A	476	29.384	61.026	72.974	1.00	17.38	6
3766	O	SER	A	476	29.103	59.902	73.358	1.00	18.38	8
3767	CB	ASER	A	476	28.931	62.557	74.900	0.70	19.66	6
3768	OG	ASER	A	476	28.330	63.803	75.257	0.70	22.12	8
3767	CB	BSER	A	476	28.870	62.465	74.932	0.30	18.16	6
3768	OG	BSER	A	476	30.220	62.822	75.127	0.30	20.85	8
3769	N	VAL	A	477	30.430	61.284	72.188	1.00	15.08	7
3770	CA	VAL	A	477	31.322	60.248	71.766	1.00	16.01	6
3771	C	VAL	A	477	32.549	60.211	72.699	1.00	14.42	6
3772	O	VAL	A	477	33.117	61.277	72.913	1.00	16.73	8
3773	CB	VAL	A	477	31.852	60.539	70.320	1.00	12.17	6
3774	CG1	VAL	A	477	32.718	59.377	69.872	1.00	15.08	6
3775	CG2	VAL	A	477	30.665	60.682	69.389	1.00	13.12	6
3776	N	ALA	A	478	32.802	59.035	73.236	1.00	17.39	6
3777	CA	ALA	A	478	33.956	58.962	74.181	1.00	17.45	6
3778	C	ALA	A	478	35.248	59.401	73.501	1.00	19.79	6
3779	O	ALA	A	478	35.398	59.204	72.298	1.00	15.95	8
3780	CB	ALA	A	478	34.031	57.543	74.731	1.00	18.76	6
3781	N	SER	A	479	36.152	60.069	74.181	1.00	15.72	7
3782	CA	SER	A	479	37.405	60.518	73.501	1.00	17.20	6
3783	C	SER	A	479	38.170	59.392	72.855	1.00	16.18	6
3784	O	SER	A	479	38.177	58.231	73.329	1.00	16.38	8
3785	CB	SER	A	479	38.262	61.218	74.569	1.00	23.12	6
3786	OG	SER	A	479	37.554	62.348	75.047	1.00	22.40	8
3787	N	PHE	A	480	38.774	59.656	71.674	1.00	16.00	7
3788	CA	PHE	A	480	39.449	58.638	70.881	1.00	15.01	6
3789	C	PHE	A	480	40.575	59.308	70.092	1.00	14.53	6
3790	O	PHE	A	480	40.568	60.548	70.019	1.00	17.08	8
3791	CB	PHE	A	480	38.538	57.834	69.938	1.00	14.92	6
3792	CG	PHE	A	480	37.888	58.695	68.846	1.00	15.98	6
3793	CD1	PHE	A	480	36.748	59.437	69.087	1.00	14.48	6
3794	CD2	PHE	A	480	38.493	58.728	67.589	1.00	14.19	8
3795	CE1	PHE	A	480	36.160	60.227	68.092	1.00	18.16	6
3796	CE2	PHE	A	480	37.899	59.513	66.609	1.00	18.21	6
3797	CZ	PHE	A	480	36.770	60.255	66.848	1.00	19.62	6
3798	N	THR	A	481	41.459	58.476	69.565	1.00	15.04	7
3799	CA	THR	A	481	42.525	59.036	68.696	1.00	12.66	6
3800	C	THR	A	481	42.186	58.783	67.223	1.00	14.82	6
3801	O	THR	A	481	41.811	57.641	66.885	1.00	16.40	6
3802	CB	THR	A	481	43.859	58.332	69.002	1.00	19.94	6
3803	OG1	THR	A	481	44.253	58.543	70.360	1.00	21.84	8

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3804  CG2 THR A 481   44.953  58.894  68.091  1.00  20.55   6
3805  N   LEU A 482   42.216  59.860  66.454  1.00  14.98   7
3806  CA  LEU A 482   42.004  59.690  64.991  1.00  13.06   6
3807  C   LEU A 482   43.383  59.527  64.394  1.00  14.32   6
3808  O   LEU A 482   44.199  60.432  64.487  1.00  15.20   8
3809  CB  LEU A 482   41.231  60.917  64.486  1.00  13.14   6
3810  CG  LEU A 482   40.812  60.936  62.999  1.00  14.96   6
3811  CD1 LEU A 482   39.938  59.717  62.720  1.00  16.24   6
3812  CD2 LEU A 482   40.073  62.242  62.727  1.00  14.36   6
3813  N   ALA A 483   43.621  58.414  63.680  1.00  12.65   6
3814  CA  ALA A 483   44.961  58.090  63.175  1.00  13.54   7
3815  C   ALA A 483   45.370  58.961  62.000  1.00  13.15   6
3816  O   ALA A 483   44.562  59.723  61.387  1.00  13.10   8
3817  CB  ALA A 483   44.892  56.581  62.782  1.00  12.82   6
3818  N   PRO A 484   46.637  58.980  61.687  1.00  14.66   7
3819  CA  PRO A 434   47.183  59.827  60.632  1.00  11.27   6
3820  C   PRO A 484   46.509  59.496  59.301  1.00  12.98   6
3821  O   PRO A 484   46.374  58.331  58.884  1.00  13.40   8
3822  CB  PRO A 484   48.704  59.546  60.545  1.00  15.48   6
3823  CG  PRO A 484   48.959  58.945  61.936  1.00  15.92   6
3824  CD  PRO A 484   47.698  58.160  62.333  1.00  15.61   6
3825  N   GLY A 485   45.977  60.548  58.671  1.00  14.03   6
3826  CA  GLY A 485   45.399  60.352  57.309  1.00  13.19   6
3827  C   GLY A 485   44.067  59.600  57.376  1.00  16.21   6
3828  O   GLY A 485   43.561  59.235  56.271  1.00  13.21   8
3829  N   ALA A 486   43.508  59.349  58.553  1.00  11.65   7
3830  CA  ALA A 486   42.375  58.395  58.545  1.00  13.98   6
3831  C   ALA A 486   41.050  58.969  58.068  1.00  15.41   6
3832  O   ALA A 486   40.730  60.148  58.196  1.00  13.33   8
3833  CB  ALA A 486   42.233  57.937  60.002  1.00  14.43   6
3834  N   VAL A 487   40.213  58.008  57.647  1.00  12.17   6
3835  CA  VAL A 487   38.776  58.220  57.403  1.00   9.79   6
3836  C   VAL A 487   38.057  57.174  58.257  1.00   9.96   6
3837  O   VAL A 487   38.406  55.970  58.179  1.00  12.30   6
3838  CB  VAL A 487   38.428  58.022  55.919  1.00  10.95   8
3839  CG1 VAL A 487   36.871  58.150  55.762  1.00  12.89   6
3840  CG2 VAL A 487   39.127  58.990  54.964  1.00  15.13   6
3841  N   SER A 488   37.112  57.635  59.078  1.00  11.84   6
3842  CA  SER A 488   36.335  56.623  59.865  1.00  12.18   6
3843  C   SER A 488   34.867  56.972  59.699  1.00  11.64   6
3844  O   SER A 488   34.519  58.154  59.650  1.00  12.81   8
3845  CB  SER A 488   36.850  56.705  61.336  1.00  15.82   6
3846  OG  SER A 488   36.431  55.439  61.907  1.00  19.75   8
3847  N   VAL A 489   33.950  55.975  59.626  1.00  11.85   7
3848  CA  VAL A 489   32.567  56.313  59.340  1.00  11.74   6
3849  C   VAL A 489   31.685  55.457  60.269  1.00  12.51   6
3850  O   VAL A 489   31.910  54.276  60.410  1.00  13.59   8
3851  CB  VAL A 489   32.231  55.903  57.882  1.00  12.50   6
3852  CG1 VAL A 489   30.712  56.154  57.650  1.00  11.14   6
3853  CG2 VAL A 489   32.978  56.812  56.882  1.00  13.14   7
3854  N   TRP A 490   30.771  56.135  60.914  1.00  11.35   6
3855  CA  TRP A 490   29.900  55.413  61.883  1.00  11.90   6
3856  C   TRP A 490   28.446  55.715  61.508  1.00  13.16   8
3857  O   TRP A 490   28.103  56.878  61.324  1.00  16.40   6
3858  CB  TRP A 490   30.222  55.953  63.283  1.00  12.61   6
3859  CG  TRP A 490   31.708  55.807  63.639  1.00  12.91   6
3860  CD1 TRP A 490   32.381  54.628  63.904  1.00  14.71   6
3861  CD2 TRP A 490   32.632  56.884  63.716  1.00  14.18   7
3862  NE1 TRP A 490   33.717  54.969  64.166  1.00  13.76   6
3863  CE2 TRP A 490   33.883  56.328  64.049  1.00  14.41   6
3864  CE3 TRP A 490   32.509  56.271  63.546  1.00  16.82   6
3865  CZ2 TRP A 490   35.020  57.117  64.216  1.00  13.91   6
3866  CZ3 TRP A 490   33.641  59.052  63.701  1.00  14.61   7
3867  CH2 TRP A 490   34.894  58.461  64.044  1.00  14.54   6
3868  N   GLN A 491   27.680  54.618  61.394  1.00  11.75   6
3869  CA  GLN A 491   26.332  54.869  60.903  1.00  12.20   6
3870  C   GLN A 491   25.221  54.121  61.637  1.00  13.63   8
3871  O   GLN A 491   25.472  53.092  62.207  1.00  13.45   7
3872  CB  GLN A 491   26.259  54.452  59.410  1.00  12.17   6
3873  CG  GLN A 491   26.541  52.972  59.117  1.00  12.78   6
3874  CD  GLN A 491   25.208  52.219  58.993  1.00  19.79   8
3875  OE1 GLN A 491   24.222  52.726  58.485  1.00  16.61   6
3876  NE2 GLN A 491   25.214  50.952  59.455  1.00  19.28   6
3877  N   TYR A 492   24.024  54.698  61.523  1.00  13.67   6
3878  CA  TYR A 492   22.819  54.043  62.092  1.00  14.21   6
3879  C   TYR A 492   21.665  54.203  61.102  1.00  17.30   8
3880  O   TYR A 492   21.507  55.341  60.614  1.00  14.76   6
3881  CB  TYR A 492   22.452  54.774  63.393  1.00  15.27   6
3882  CG  TYR A 492   21.152  54.219  63.984  1.00  15.56   6
3883  CD1 TYR A 492   21.146  53.010  64.638  1.00  20.95   6
3884  CD2 TYR A 492   19.963  54.959  63.800  1.00  18.48   6
3885  CE1 TYR A 492   19.956  52.511  65.176  1.00  21.55   6
3886  CE2 TYR A 492   18.770  54.444  64.324  1.00  19.51   6
3887  CZ  TYR A 492   18.817  53.254  65.003  1.00  22.52   6
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Serial	Atom	Residue	Chain	ResSeq	X	Y	Z	Occ	B	Elem
3888	OH	TYR	A	492	17.596	52.781	65.511	1.00	24.89	8
3889	N	SER	A	493	20.879	53.151	60.925	1.00	15.17	7
3890	CA	SER	A	493	19.666	53.363	60.096	1.00	14.91	6
3891	C	SER	A	493	18.548	52.462	60.616	1.00	20.17	6
3892	O	SER	A	493	18.887	51.457	61.265	1.00	22.90	8
3893	CB	SER	A	493	19.867	53.083	58.619	1.00	23.97	6
3894	OG	SER	A	493	20.148	51.710	58.532	1.00	30.62	8
3895	N	THR	A	494	17.344	52.991	60.393	1.00	15.89	7
3896	CA	THR	A	494	16.243	52.127	60.906	1.00	18.40	6
3897	C	THR	A	494	14.999	52.406	60.060	1.00	20.14	6
3898	O	THR	A	494	14.967	53.429	59.367	1.00	20.45	8
3899	CB	THR	A	494	15.993	52.424	62.396	1.00	24.38	6
3900	OG1	THR	A	494	15.137	51.363	62.864	1.00	25.66	8
3901	CG2	THR	A	494	15.368	53.751	62.737	1.00	27.09	6
3902	N	SER	A	495	14.017	51.534	60.157	1.00	22.56	7
3903	CA	SER	A	495	12.772	51.870	59.422	1.00	19.72	6
3904	C	SER	A	495	12.082	53.049	60.021	1.00	19.50	6
3905	O	SER	A	495	12.132	53.418	61.188	1.00	21.11	8
3906	CB	ASER	A	495	11.766	50.698	59.444	0.60	26.59	6
3907	OG	ASER	A	495	12.447	49.487	59.259	0.60	31.40	8
3906	CB	BSER	A	495	11.888	50.603	59.441	0.40	20.10	6
3907	OG	BSER	A	495	11.922	50.184	60.798	0.40	20.04	8
3908	N	ALA	A	496	11.315	53.727	59.141	1.00	20.53	7
3909	CA	ALA	A	496	10.529	54.893	59.493	1.00	21.55	6
3910	C	ALA	A	496	9.094	54.492	59.960	1.00	24.11	6
3911	O	ALA	A	496	8.536	53.596	59.350	1.00	29.93	8
3912	CB	ALA	A	496	10.354	55.717	58.189	1.00	22.92	6
3913	N	SER	A	497	8.599	55.279	60.902	1.00	26.84	7
3914	CA	SER	A	497	7.226	54.967	61.354	1.00	32.64	6
3915	C	SER	A	497	6.242	56.009	60.853	1.00	31.20	6
3916	O	SER	A	497	5.049	55.788	61.090	1.00	33.40	8
3917	CB	SER	A	497	7.183	54.836	62.875	1.00	35.13	6
3918	OG	SER	A	497	7.578	56.030	63.515	1.00	39.06	8
3919	N	ALA	A	498	6.685	56.920	59.967	1.00	25.21	7
3920	CA	ALA	A	498	5.749	57.878	59.354	1.00	21.05	6
3921	C	ALA	A	498	6.350	58.189	57.975	1.00	22.00	6
3922	O	ALA	A	498	7.541	57.907	57.763	1.00	18.25	8
3923	CB	ALA	A	498	5.737	59.103	60.231	1.00	22.15	6
3924	N	PRO	A	499	5.633	58.771	57.054	1.00	18.84	7
3925	CA	PRO	A	499	6.176	59.058	55.718	1.00	19.31	6
3926	C	PRO	A	499	7.174	60.210	55.762	1.00	15.69	6
3927	O	PRO	A	499	6.934	61.242	56.370	1.00	16.96	8
3928	CB	PRO	A	499	4.930	59.294	54.823	1.00	20.07	6
3929	CG	PRO	A	499	3.927	59.790	55.829	1.00	23.72	6
3930	CD	PRO	A	499	4.205	59.113	57.173	1.00	23.70	6
3931	N	GLN	A	500	8.288	59.975	55.037	1.00	18.24	7
3932	CA	GLN	A	500	9.423	60.939	55.075	1.00	13.65	6
3933	C	GLN	A	500	9.921	61.084	53.596	1.00	12.93	6
3934	O	GLN	A	500	10.256	60.065	53.014	1.00	15.78	8
3935	CB	GLN	A	500	10.601	60.286	55.860	1.00	15.18	6
3936	CG	GLN	A	500	10.189	60.048	57.328	1.00	16.32	6
3937	CD	GLN	A	500	11.284	59.264	58.126	1.00	15.95	6
3938	OE1	GLN	A	500	12.239	58.781	57.571	1.00	17.99	8
3939	NE2	GLN	A	500	11.008	59.238	59.419	1.00	20.52	7
3940	N	ILE	A	501	9.662	62.283	53.054	1.00	12.81	7
3941	CA	ILE	A	501	10.101	62.501	51.667	1.00	12.56	6
3942	C	ILE	A	501	11.594	62.914	51.642	1.00	13.47	6
3943	O	ILE	A	501	11.899	63.905	52.301	1.00	14.35	8
3944	CB	ILE	A	501	9.262	63.635	51.032	1.00	12.51	6
3945	CG1	ILE	A	501	7.788	63.216	50.842	1.00	14.81	6
3946	CG2	ILE	A	501	9.888	63.939	49.629	1.00	12.71	6
3947	CD1	ILE	A	501	6.897	64.308	50.244	1.00	15.70	6
3948	N	GLY	A	502	12.383	62.169	50.872	1.00	12.33	7
3949	CA	GLY	A	502	13.793	62.673	50.746	1.00	12.41	6
3950	C	GLY	A	502	13.915	63.416	49.403	1.00	11.68	6
3951	O	GLY	A	502	14.806	64.300	49.332	1.00	12.18	8
3952	N	SER	A	503	13.120	63.068	48.395	1.00	12.99	7
3953	CA	SER	A	503	13.334	63.710	47.073	1.00	12.73	6
3954	C	SER	A	503	11.970	63.672	46.342	1.00	14.01	6
3955	O	SER	A	503	11.263	62.678	46.449	1.00	14.22	8
3956	CB	SER	A	503	14.364	62.853	46.296	1.00	12.66	6
3957	OG	SER	A	503	14.467	63.267	44.940	1.00	12.94	8
3958	N	VAL	A	504	11.736	64.763	45.625	1.00	12.13	7
3959	CA	VAL	A	504	10.738	64.716	44.521	1.00	11.27	6
3960	C	VAL	A	504	11.521	65.159	43.243	1.00	11.68	6
3961	O	VAL	A	504	12.287	66.121	43.321	1.00	11.43	8
3962	CB	VAL	A	504	9.614	65.700	44.790	1.00	12.52	6
3963	CG1	VAL	A	504	8.670	65.766	43.563	1.00	15.49	6
3964	CG2	VAL	A	504	8.740	65.263	46.021	1.00	12.38	6
3965	N	ALA	A	505	11.424	64.285	42.236	1.00	11.00	7
3966	CA	ALA	A	505	12.174	64.595	41.003	1.00	10.56	6
3967	C	ALA	A	505	11.447	64.174	39.798	1.00	12.28	6
3968	O	ALA	A	505	10.738	63.165	39.919	1.00	13.54	8
3969	CB	ALA	A	505	13.567	63.894	41.072	1.00	11.90	6

Atom	Name	Res	Chain	Seq	X	Y	Z	Occ	B	El
3970	N	PRO	A	506	11.674	64.766	38.661	1.00	12.95	7
3971	CA	PRO	A	506	12.353	66.020	38.410	1.00	11.12	6
3972	C	PRO	A	506	11.747	67.182	39.194	1.00	12.55	6
3973	O	PRO	A	506	10.711	67.015	39.862	1.00	12.33	8
3974	CB	PRO	A	506	12.227	66.313	36.887	1.00	13.30	6
3975	CG	PRO	A	506	11.545	65.090	36.310	1.00	15.76	6
3976	CD	PRO	A	506	11.007	64.329	37.471	1.00	14.19	6
3977	N	ASN	A	507	12.364	68.381	39.142	1.00	10.92	7
3978	CA	ASN	A	507	11.835	69.493	39.909	1.00	11.12	6
3979	C	ASN	A	507	10.940	70.447	39.087	1.00	10.54	6
3980	O	ASN	A	507	10.497	71.416	39.701	1.00	12.30	8
3981	CB	ASN	A	507	13.043	70.353	40.401	1.00	11.79	6
3982	CG	ASN	A	507	14.033	69.459	41.196	1.00	14.07	6
3983	OD1	ASN	A	507	15.192	69.361	40.778	1.00	13.37	8
3984	ND2	ASN	A	507	13.592	68.845	42.261	1.00	13.47	7
3985	N	MET	A	508	10.654	70.046	37.850	1.00	10.20	7
3986	CA	MET	A	508	9.823	70.957	37.021	1.00	11.04	6
3987	C	MET	A	508	9.102	70.087	36.027	1.00	11.63	6
3988	O	MET	A	508	9.633	69.029	35.633	1.00	12.66	8
3989	CB	MET	A	508	10.929	71.782	36.219	1.00	12.80	6
3990	CG	MET	A	508	10.270	72.808	35.305	1.00	15.57	6
3991	SD	MET	A	508	11.558	73.692	34.325	1.00	12.60	16
3992	CE	MET	A	508	11.921	72.582	33.003	1.00	11.82	6
3993	N	GLY	A	509	7.935	70.545	35.550	1.00	11.36	7
3994	CA	GLY	A	509	7.253	69.739	34.528	1.00	10.02	6
3995	C	GLY	A	509	5.851	70.331	34.276	1.00	13.18	6
3996	O	GLY	A	509	5.506	71.321	34.889	1.00	13.82	8
3997	N	ILE	A	510	5.070	69.599	33.480	1.00	11.23	7
3998	CA	ILE	A	510	3.674	70.061	33.192	1.00	12.19	6
3999	C	ILE	A	510	2.746	68.964	33.701	1.00	11.61	6
4000	O	ILE	A	510	3.156	67.806	33.903	1.00	13.58	8
4001	CB	ILE	A	510	3.456	70.272	31.683	1.00	12.75	6
4002	CG1	ILE	A	510	3.840	69.011	30.868	1.00	11.82	6
4003	CG2	ILE	A	510	4.241	71.485	31.163	1.00	14.82	6
4004	CD1	ILE	A	510	3.293	69.108	29.417	1.00	18.53	6
4005	N	PRO	A	511	1.462	69.267	33.824	1.00	13.51	7
4006	CA	PRO	A	511	0.442	68.291	34.164	1.00	14.15	6
4007	C	PRO	A	511	0.531	67.082	33.263	1.00	13.29	6
4008	O	PRO	A	511	0.780	67.180	32.026	1.00	13.86	8
4009	CB	PRO	A	511	-0.913	69.037	33.810	1.00	13.10	6
4010	CG	PRO	A	511	-0.528	70.435	34.265	1.00	15.94	6
4011	CD	PRO	A	511	0.925	70.625	33.729	1.00	16.03	6
4012	N	GLY	A	512	0.462	65.892	33.876	1.00	12.55	7
4013	CA	GLY	A	512	0.554	64.629	33.165	1.00	14.21	6
4014	C	GLY	A	512	1.950	63.971	33.170	1.00	13.76	6
4015	O	GLY	A	512	2.002	62.765	32.923	1.00	15.31	8
4016	N	ASN	A	513	2.994	64.789	33.454	1.00	12.18	7
4017	CA	ASN	A	513	4.306	64.154	33.561	1.00	10.79	6
4018	C	ASN	A	513	4.360	63.144	34.731	1.00	15.27	6
4019	O	ASN	A	513	3.699	63.392	35.738	1.00	13.64	8
4020	CB	ASN	A	513	5.408	65.230	33.833	1.00	11.43	6
4021	CG	ASN	A	513	5.762	66.023	32.583	1.00	11.94	6
4022	OD1	ASN	A	513	6.738	66.874	32.746	1.00	16.21	8
4023	ND2	ASN	A	513	5.078	65.931	31.505	1.00	8.90	7
4024	N	VAL	A	514	5.280	62.177	34.576	1.00	11.74	7
4025	CA	VAL	A	514	5.505	61.269	35.745	1.00	11.63	6
4026	C	VAL	A	514	6.594	61.894	36.612	1.00	13.73	6
4027	O	VAL	A	514	7.617	62.379	36.131	1.00	15.35	8
4028	CB	VAL	A	514	6.072	59.931	35.221	1.00	13.67	6
4029	CG1	VAL	A	514	6.529	59.036	36.390	1.00	14.35	6
4030	CG2	VAL	A	514	5.042	59.176	34.393	1.00	16.73	6
4031	N	VAL	A	515	6.335	61.959	37.923	1.00	10.41	7
4032	CA	VAL	A	515	7.256	62.496	38.927	1.00	12.67	6
4033	C	VAL	A	515	7.433	61.442	40.001	1.00	13.31	6
4034	O	VAL	A	515	6.495	60.711	40.306	1.00	14.78	8
4035	CB	VAL	A	515	6.563	63.772	39.384	1.00	14.30	6
4036	CG1	VAL	A	515	7.228	64.271	40.775	1.00	18.11	6
4037	CG2	VAL	A	515	6.678	64.883	38.435	1.00	16.90	6
4038	N	THR	A	516	8.669	61.321	40.514	1.00	11.19	7
4039	CA	THR	A	516	8.900	60.252	41.495	1.00	11.78	6
4040	C	THR	A	516	9.271	60.860	42.835	1.00	11.71	6
4041	O	THR	A	516	10.092	61.763	42.959	1.00	12.89	8
4042	CB	THR	A	516	10.107	59.384	41.001	1.00	14.88	6
4043	OG1	THR	A	516	9.696	58.789	39.742	1.00	14.59	8
4044	CG2	THR	A	516	10.446	58.275	42.004	1.00	14.18	6
4045	N	ILE	A	517	8.600	60.281	43.863	1.00	9.66	7
4046	CA	ILE	A	517	8.878	60.681	45.251	1.00	10.17	6
4047	C	ILE	A	517	9.670	59.586	45.920	1.00	13.57	6
4048	O	ILE	A	517	9.172	58.427	45.949	1.00	13.82	8
4049	CB	ILE	A	517	7.493	60.899	45.948	1.00	12.16	6
4050	CG1	ILE	A	517	6.659	61.984	45.278	1.00	15.38	6
4051	CG2	ILE	A	517	7.745	61.419	47.392	1.00	14.69	6
4052	CD1	ILE	A	517	5.200	61.951	45.779	1.00	22.16	6
4053	N	ASP	A	518	10.911	59.861	46.363	1.00	13.29	7

#	Atom	Res	Chain	Seq	X	Y	Z	Occ	B	n
4054	CA	ASP	A	518	11.743	58.802	46.976	1.00	14.12	6
4055	C	ASP	A	518	11.868	59.161	48.435	1.00	13.11	6
4056	O	ASP	A	518	11.922	60.347	48.791	1.00	14.37	8
4057	CB	ASP	A	518	13.159	58.791	46.351	1.00	14.49	6
4058	CG	ASP	A	518	13.106	58.207	44.962	1.00	17.62	6
4059	OD1	ASP	A	518	12.858	56.963	44.888	1.00	17.47	8
4060	OD2	ASP	A	518	13.289	58.874	43.931	1.00	16.96	8
4061	N	GLY	A	519	11.769	58.138	49.315	1.00	13.39	7
4062	CA	GLY	A	519	11.872	58.484	50.757	1.00	14.21	6
4063	C	GLY	A	519	11.716	57.152	51.549	1.00	13.61	6
4064	O	GLY	A	519	12.278	56.130	51.134	1.00	15.70	8
4065	N	LYS	A	520	11.059	57.319	52.696	1.00	17.22	7
4066	CA	LYS	A	520	10.867	56.091	53.545	1.00	15.85	6
4067	C	LYS	A	520	9.539	56.230	54.256	1.00	16.54	6
4068	O	LYS	A	520	9.004	57.334	54.385	1.00	17.43	8
4069	CB	LYS	A	520	11.866	56.344	54.747	1.00	17.75	6
4070	CG	LYS	A	520	13.318	56.307	54.421	1.00	24.98	6
4071	CD	LYS	A	520	14.147	56.512	55.698	1.00	22.21	6
4072	CE	LYS	A	520	14.074	55.145	56.459	1.00	23.54	6
4073	NZ	LYS	A	520	15.426	54.879	57.005	1.00	27.08	7
4074	N	GLY	A	521	9.021	55.065	54.728	1.00	16.42	7
4075	CA	GLY	A	521	7.870	55.223	55.627	1.00	16.98	6
4076	C	GLY	A	521	6.540	55.347	54.874	1.00	19.49	6
4077	O	GLY	A	521	5.533	55.716	55.525	1.00	18.22	8
4078	N	PHE	A	522	6.569	55.053	53.560	1.00	19.15	7
4079	CA	PHE	A	522	5.299	55.230	52.839	1.00	17.53	6
4080	C	PHE	A	522	4.344	54.037	53.029	1.00	20.57	6
4081	O	PHE	A	522	3.173	54.222	52.638	1.00	20.45	8
4082	CB	PHE	A	522	5.587	55.448	51.342	1.00	17.55	6
4083	CG	PHE	A	522	6.513	56.605	51.009	1.00	15.40	6
4084	CD1	PHE	A	522	6.601	57.740	51.763	1.00	17.48	6
4085	CD2	PHE	A	522	7.262	56.452	49.824	1.00	17.80	6
4086	CE1	PHE	A	522	7.480	58.778	51.376	1.00	20.21	6
4087	CE2	PHE	A	522	8.142	57.475	49.422	1.00	16.35	6
4088	CZ	PHE	A	522	8.257	58.607	50.214	1.00	14.66	7
4089	N	GLY	A	523	4.826	52.902	53.488	1.00	23.38	6
4090	CA	GLY	A	523	3.947	51.721	53.586	1.00	26.21	6
4091	C	GLY	A	523	3.753	51.037	52.260	1.00	27.39	6
4092	O	GLY	A	523	4.082	51.536	51.167	1.00	21.64	8
4093	N	THR	A	524	3.204	49.785	52.316	1.00	26.00	7
4094	CA	THR	A	524	3.043	49.010	51.103	1.00	24.30	6
4095	C	THR	A	524	1.636	49.241	50.556	1.00	26.74	6
4096	O	THR	A	524	1.433	48.984	49.372	1.00	34.03	8
4097	CB	THR	A	524	3.314	47.505	51.308	1.00	36.63	6
4098	OG1	THR	A	524	2.403	47.020	52.315	1.00	34.71	8
4099	CG2	THR	A	524	4.711	47.251	51.862	1.00	34.40	6
4100	N	THR	A	525	0.748	49.673	51.432	1.00	27.38	7
4101	CA	THR	A	525	-0.646	49.948	51.027	1.00	29.73	6
4102	C	THR	A	525	-0.696	51.367	50.456	1.00	29.44	6
4103	O	THR	A	525	-0.315	52.313	51.159	1.00	28.46	8
4104	CB	THR	A	525	-1.558	49.835	52.265	1.00	31.59	6
4105	OG1	THR	A	525	-1.416	48.485	52.779	1.00	34.50	8
4106	CG2	THR	A	525	-3.020	50.076	51.920	1.00	33.53	6
4107	N	GLN	A	526	-1.341	51.530	49.304	1.00	26.38	7
4108	CA	GLN	A	526	-1.383	52.864	48.695	1.00	27.25	6
4109	C	GLN	A	526	-2.090	53.867	49.574	1.00	29.32	6
4110	O	GLN	A	526	-3.264	53.665	49.960	1.00	25.61	8
4111	CB	GLN	A	526	-2.101	52.757	47.340	1.00	30.46	6
4112	CG	GLN	A	526	-2.028	53.996	46.486	1.00	28.17	6
4113	CD	GLN	A	526	-2.542	53.744	45.055	1.00	30.46	6
4114	OE1	GLN	A	526	-3.419	54.518	44.679	1.00	28.73	8
4115	NE2	GLN	A	526	-1.951	52.750	44.438	1.00	33.21	7
4116	N	GLY	A	527	-1.476	55.032	49.820	1.00	31.17	7
4117	CA	GLY	A	527	-2.091	56.150	49.471	1.00	22.05	6
4118	C	GLY	A	527	-2.415	57.258	48.405	1.00	20.00	6
4119	O	GLY	A	527	-2.894	56.897	49.804	1.00	19.58	8
4120	N	THR	A	528	-2.136	58.506	48.884	1.00	21.95	7
4121	CA	THR	A	528	-2.428	59.597	48.808	1.00	21.30	6
4122	C	THR	A	528	-1.268	60.584	49.647	1.00	18.42	6
4123	O	THR	A	528	-0.375	60.642	49.320	1.00	21.48	8
4124	CB	THR	A	528	-3.685	60.387	50.657	1.00	20.31	6
4125	OG1	THR	A	528	-3.522	60.782	49.240	1.00	24.98	8
4126	CG2	THR	A	528	-4.933	59.511	50.657	1.00	30.73	6
4127	N	VAL	A	529	-1.209	61.255	47.667	1.00	26.25	7
4128	CA	VAL	A	529	-0.167	62.275	47.425	1.00	16.67	6
4129	C	VAL	A	529	-0.924	63.532	46.987	1.00	15.19	6
4130	O	VAL	A	529	-1.825	63.328	46.155	1.00	16.18	8
4131	CB	VAL	A	529	0.700	61.851	46.216	1.00	16.64	6
4132	CG1	VAL	A	529	1.664	62.995	45.841	1.00	14.16	6
4133	CG2	VAL	A	529	1.483	60.594	46.604	1.00	17.46	6
4134	N	THR	A	530	-0.533	64.669	47.531	1.00	18.23	7
4135	CA	THR	A	530	-1.148	65.900	47.016	1.00	14.50	6
4136	C	THR	A	530	-0.038	66.814	46.492	1.00	14.78	6
4137	O	THR	A	530	1.076	66.792	47.027	1.00	17.57	8

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4138  CB   THR A 530   -1.954  66.621  48.090  1.00  15.93  6
4139  OG1  THR A 530   -1.209  66.722  49.308  1.00  17.69  8
4140  CG2  THR A 530   -3.279  65.858  48.340  1.00  17.39  6
4141  N    PHE A 531   -0.395  67.691  45.569  1.00  14.04  7
4142  CA   PHE A 531    0.458  68.822  45.142  1.00  13.01  6
4143  C    PHE A 531   -0.344  70.073  45.496  1.00  15.03  6
4144  O    PHE A 531   -1.454  70.292  44.989  1.00  16.16  8
4145  CB   PHE A 531    0.659  68.823  43.604  1.00  13.87  6
4146  CG   PHE A 531    1.611  67.777  43.040  1.00  13.21  6
4147  CD1  PHE A 531    1.438  66.431  43.191  1.00  14.34  6
4148  CD2  PHE A 531    2.662  68.228  42.240  1.00  14.96  6
4149  CE1  PHE A 531    2.288  65.515  42.629  1.00  16.70  6
4150  CE2  PHE A 531    3.545  67.306  41.691  1.00  13.81  6
4151  CZ   PHE A 531    3.385  65.943  41.836  1.00  16.70  6
4152  N    GLY A 532    0.118  70.806  46.490  1.00  14.18  7
4153  CA   GLY A 532   -0.569  72.077  46.884  1.00  15.57  6
4154  C    GLY A 532   -1.992  71.702  47.378  1.00  19.88  6
4155  O    GLY A 532   -2.928  72.482  47.068  1.00  18.91  8
4156  N    GLY A 533   -2.193  70.510  47.921  1.00  17.41  7
4157  CA   GLY A 533   -3.524  70.089  48.400  1.00  18.31  6
4158  C    GLY A 533   -4.368  69.372  47.370  1.00  20.08  6
4159  O    GLY A 533   -5.463  68.817  47.637  1.00  19.31  8
4160  N    VAL A 534   -3.923  69.391  46.097  1.00  16.06  7
4161  CA   VAL A 534   -4.592  68.721  44.999  1.00  15.11  6
4162  C    VAL A 534   -4.197  67.275  44.894  1.00  15.67  6
4163  O    VAL A 534   -3.019  66.888  44.712  1.00  17.94  8
4164  CB   VAL A 534   -4.368  69.480  43.645  1.00  13.67  6
4165  CG1  VAL A 534   -5.101  68.739  42.509  1.00  15.45  6
4166  CG2  VAL A 534   -4.807  70.937  43.768  1.00  16.93  6
4167  N    THR A 535   -5.185  66.334  44.967  1.00  13.93  7
4168  CA   THR A 535   -4.827  64.927  44.890  1.00  16.87  6
4169  C    THR A 535   -4.271  64.521  43.536  1.00  20.28  6
4170  O    THR A 535   -4.796  64.893  42.462  1.00  17.64  8
4171  CB   THR A 535   -6.065  64.042  45.192  1.00  20.23  6
4172  OG1  THR A 535   -6.446  64.284  46.576  1.00  20.94  8
4173  CG2  THR A 535   -5.787  62.565  45.026  1.00  24.73  6
4174  N    ALA A 536   -3.162  63.773  43.562  1.00  15.87  7
4175  CA   ALA A 536   -2.521  63.324  42.337  1.00  17.27  6
4176  C    ALA A 536   -2.808  61.859  42.042  1.00  20.29  6
4177  O    ALA A 536   -2.929  61.111  43.030  1.00  20.32  8
4178  CB   ALA A 536   -0.976  63.441  42.384  1.00  17.03  6
4179  N    THR A 537   -2.937  61.461  40.791  1.00  16.03  7

4180  CA   THR A 537   -3.151  60.043  40.489  1.00  15.38  6
4181  C    THR A 537   -1.865  59.242  40.679  1.00  19.21  6
4182  O    THR A 537   -0.823  59.697  40.166  1.00  19.02  8
4183  CB   THR A 537   -3.564  59.955  38.998  1.00  18.90  6
4184  OG1  THR A 537   -4.828  60.651  38.950  1.00  19.30  8
4185  CG2  THR A 537   -3.697  58.508  38.591  1.00  18.81  6
4186  N    VAL A 538   -1.925  58.176  41.415  1.00  16.51  7
4187  CA   VAL A 538   -0.704  57.394  41.746  1.00  16.38  6
4188  C    VAL A 538   -0.516  56.415  40.613  1.00  18.34  6
4189  O    VAL A 538   -1.390  55.599  40.252  1.00  20.80  8
4190  CB   VAL A 538   -0.896  56.665  43.080  1.00  17.97  6
4191  CG1  VAL A 538    0.219  55.621  43.337  1.00  16.03  6
4192  CG2  VAL A 538   -1.016  57.646  44.226  1.00  19.59  6
4193  N    LYS A 539    0.696  56.340  40.055  1.00  15.64  7
4194  CA   LYS A 539    1.119  55.341  39.108  1.00  14.97  6
4195  C    LYS A 539    1.626  54.044  39.732  1.00  17.57  6
4196  O    LYS A 539    1.313  52.885  39.375  1.00  18.05  8
4197  CB   LYS A 539    2.264  55.914  38.209  1.00  17.15  6
4198  CG   LYS A 539    2.814  54.859  37.246  1.00  20.63  6
4199  CD   LYS A 539    3.860  55.636  36.368  1.00  25.39  6
4200  CE   LYS A 539    3.601  55.199  34.949  1.00  41.98  6
4201  NZ   LYS A 539    4.369  53.976  34.672  1.00  30.47  7
4202  N    SER A 540    2.424  54.212  40.787  1.00  15.53  7
4203  CA   SER A 540    2.919  53.073  41.587  1.00  16.52  6
4204  C    SER A 540    3.231  53.502  42.999  1.00  17.69  6
4205  O    SER A 540    3.482  54.680  43.306  1.00  16.82  8
4206  CB   SER A 540    4.136  52.424  40.903  1.00  20.58  6
4207  OG   SER A 540    5.270  53.317  41.043  1.00  19.05  8
4208  N    TRP A 541    3.206  52.536  43.953  1.00  16.64  6
4209  CA   TRP A 541    3.361  52.905  45.378  1.00  15.40  6
4210  C    TRP A 541    4.148  51.785  46.053  1.00  20.16  6
4211  O    TRP A 541    3.682  50.647  46.008  1.00  21.11  8
4212  CB   TRP A 541    2.034  53.084  46.101  1.00  17.49  6
4213  CG   TRP A 541    2.124  53.605  47.502  1.00  15.27  6
4214  CD1  TRP A 541    2.645  52.924  48.584  1.00  18.64  6
4215  CD2  TRP A 541    1.689  54.854  48.006  1.00  15.30  6
4216  NE1  TRP A 541    2.542  53.673  49.715  1.00  19.85  6
4217  CE2  TRP A 541    1.976  54.894  49.381  1.00  17.27  6
4218  CE3  TRP A 541    1.086  55.984  47.440  1.00  15.92  6
4219  CZ2  TRP A 541    1.703  55.983  50.201  1.00  18.49  6
4220  CZ3  TRP A 541    0.787  57.054  48.223  1.00  19.66  6
4221  CH2  TRP A 541    1.076  57.063  49.619  1.00  21.41  6
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4222  N   THR A 542   5.297  52.107  46.615  1.00  18.77  7
4223  CA  THR A 542   6.111  51.224  47.437  1.00  20.27  6
4224  C   THR A 542   6.477  51.972  48.690  1.00  21.74  6
4225  O   THR A 542   6.369  53.228  48.814  1.00  17.27  8
4226  CB  THR A 542   7.356  50.641  46.743  1.00  24.92  6
4227  OG1 THR A 542   8.305  51.745  46.576  1.00  21.04  8
4228  CG2 THR A 542   7.091  49.930  45.442  1.00  26.69  6
4229  N   SER A 543   7.123  51.252  49.648  1.00  18.34  7
4230  CA  SER A 543   7.474  51.808  50.923  1.00  18.96  6
4231  C   SER A 543   8.463  52.975  50.734  1.00  17.37  6
4232  O   SER A 543   8.525  53.808  51.615  1.00  19.43  8
4233  CB  SER A 543   8.201  50.711  51.743  1.00  24.79  6
4234  OG  SER A 543   7.254  49.673  51.954  1.00  38.42  8
4235  N   ASN A 544   9.313  52.865  49.721  1.00  17.78  7
4236  CA  ASN A 544  10.349  53.917  49.575  1.00  14.67  6
4237  C   ASN A 544  10.208  54.723  48.287  1.00  15.24  6
4238  O   ASN A 544  11.018  55.668  48.072  1.00  16.63  8
4239  CB  AASN A 544 11.734  53.252  49.583  0.50  18.64  6
4240  CG  AASN A 544 12.145  52.868  51.005  0.50  24.93  6
4241  OD1 AASN A 544 11.394  53.024  51.976  0.50  27.01  8
4242  ND2 AASN A 544 13.359  52.364  51.118  0.50  19.54  7
4240  CB  BASN A 544 11.746  53.263  49.523  0.50  15.95  6
4241  CG  BASN A 544 11.998  52.552  50.860  0.50  20.57  6
4242  OD1 BASN A 544 12.195  53.205  51.884  0.50  23.06  8
4243  ND2 BASN A 544 11.914  51.250  50.767  0.50  19.16  7
4243  N   ARG A 545   9.224  54.409  47.426  1.00  14.40  7
4244  CA  ARG A 545   9.193  55.190  46.167  1.00  16.44  6
4245  C   ARG A 545   7.727  55.246  45.682  1.00  19.39  6
4246  O   ARG A 545   7.083  54.204  45.539  1.00  17.25  8
4247  CB  ARG A 545  10.085  54.589  45.084  1.00  17.96  6
4248  CG  ARG A 545   9.964  55.404  43.794  1.00  17.03  6
4249  CD  ARG A 545  10.778  54.728  42.653  1.00  15.55  6
4250  NE  ARG A 545  12.186  54.934  43.045  1.00  17.92  7
4251  CZ  ARG A 545  13.164  54.094  42.735  1.00  27.70  6
4252  NH1 ARG A 545  12.923  53.020  41.999  1.00  26.35  7
4253  NH2 ARG A 545  14.392  54.343  43.179  1.00  26.80  7
4254  N   ILE A 546   7.282  56.458  45.332  1.00  18.00  7
4255  CA  ILE A 546   5.908  56.613  44.793  1.00  14.23  6
4256  C   ILE A 546   6.077  57.267  43.406  1.00  17.86  6
4257  O   ILE A 546   6.771  58.296  43.389  1.00  15.62  8
4258  CB  ILE A 546   5.039  57.498  45.668  1.00  13.38  6
4259  CG1 ILE A 546   4.895  56.781  47.045  1.00  15.14  6

4260  CG2 ILE A 546   3.624  57.682  45.030  1.00  14.12  6
4261  CD1 ILE A 546   4.473  57.773  48.119  1.00  13.44  6
4262  N   GLU A 547   5.424  56.777  42.400  1.00  14.74  7
4263  CA  GLU A 547   5.338  57.543  41.134  1.00  14.58  6
4264  C   GLU A 547   3.943  58.116  40.945  1.00  15.30  6
4265  O   GLU A 547   2.977  57.405  41.270  1.00  15.50  8
4266  CB  GLU A 547   5.537  56.644  39.913  1.00  14.75  6
4267  CG  GLU A 547   6.987  56.074  39.890  1.00  17.63  6
4268  CD  GLU A 547   7.105  55.058  38.781  1.00  21.54  6
4269  OE1 GLU A 547   6.335  54.046  38.730  1.00  18.28  8
4270  OE2 GLU A 547   7.924  55.225  37.834  1.00  16.92  8
4271  N   VAL A 548   3.865  59.377  40.603  1.00  12.67  7
4272  CA  VAL A 548   2.565  60.033  40.428  1.00  13.16  6
4273  C   VAL A 548   3.587  60.764  39.095  1.00  15.52  6
4274  O   VAL A 548   3.587  61.028  38.477  1.00  14.55  8
4275  CB  VAL A 548   2.266  61.100  41.499  1.00  15.22  6
4276  CG1 VAL A 548   2.134  60.409  42.872  1.00  17.06  6
4277  CG2 VAL A 548   3.376  62.179  41.584  1.00  16.26  6
4278  N   TYR A 549   1.338  61.119  38.644  1.00  12.92  7
4279  CA  TYR A 549   1.226  61.997  37.481  1.00  14.43  6
4280  C   TYR A 549   0.902  63.397  37.975  1.00  13.16  6
4281  O   TYR A 549   0.223  63.571  39.016  1.00  14.84  8
4282  CB  TYR A 549   0.000  61.611  36.605  1.00  14.89  6
4283  CG  TYR A 549   0.208  60.240  36.037  1.00  14.21  6
4284  CD1 TYR A 549   1.049  60.058  34.934  1.00  17.04  6
4285  CD2 TYR A 549  -0.398  59.160  36.628  1.00  14.96  6
4286  CE1 TYR A 549   1.252  58.788  34.395  1.00  19.51  6
4287  CE2 TYR A 549  -0.214  57.885  36.081  1.00  20.11  6
4288  CZ  TYR A 549   0.577  57.730  34.984  1.00  20.28  6
4289  OH  TYR A 549   0.789  56.436  34.508  1.00  21.51  8
4290  N   VAL A 550   1.626  64.446  37.496  1.00  12.41  7
4291  CA  VAL A 550   1.317  65.794  37.957  1.00  13.30  6
4292  C   VAL A 550  -0.145  66.139  37.602  1.00  12.86  6
4293  O   VAL A 550  -0.589  65.893  36.503  1.00  15.42  8
4294  CB  VAL A 550   2.195  66.791  37.131  1.00  12.38  6
4295  CG1 VAL A 550   1.968  68.236  37.581  1.00  13.91  6
4296  CG2 VAL A 550   3.657  66.404  37.542  1.00  16.31  6
4297  N   PRO A 551  -0.828  66.685  38.603  1.00  13.48  7
4298  CA  PRO A 551  -2.272  66.942  38.387  1.00  16.42  6
4299  C   PRO A 551  -2.447  68.081  37.424  1.00  16.09  6
4300  O   PRO A 551  -1.599  68.955  37.236  1.00  16.43  8
4301  CB  PRO A 551  -2.869  67.337  39.755  1.00  20.53  6
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4302	CG	PRO	A	551	-1.772	66.991	40.707	1.00	20.55	6
4303	CD	PRO	A	551	-0.427	66.849	39.973	1.00	17.00	6
4304	N	ASN	A	552	-3.658	68.150	36.821	1.00	15.47	7
4305	CA	ASN	A	552	-4.017	69.248	35.941	1.00	15.18	7
4306	C	ASN	A	552	-4.401	70.475	36.748	1.00	19.54	6
4307	O	ASN	A	552	-5.630	70.744	36.917	1.00	19.65	8
4308	CB	ASN	A	552	-5.198	68.759	35.075	1.00	19.00	6
4309	CG	ASN	A	552	-5.522	69.706	33.925	1.00	23.61	6
4310	OD1	ASN	A	552	-4.763	70.583	33.553	1.00	29.14	8
4311	ND2	ASN	A	552	-6.635	69.481	33.239	1.00	23.65	7
4312	N	MET	A	553	-3.487	71.146	37.402	1.00	13.46	6
4313	CA	MET	A	553	-3.722	72.208	38.346	1.00	11.50	6
4314	C	MET	A	553	-3.003	73.456	37.901	1.00	15.54	6
4315	O	MET	A	553	-2.319	73.427	36.881	1.00	17.61	8
4316	CB	MET	A	553	-3.328	71.835	39.803	1.00	16.46	6
4317	CG	MET	A	553	-1.826	71.490	39.883	1.00	14.92	6
4318	SD	MET	A	553	-1.364	70.962	41.579	1.00	17.71	16
4319	CE	MET	A	553	-1.416	72.450	42.426	1.00	16.59	6
4320	N	ALA	A	554	-3.278	74.532	38.619	1.00	17.19	7
4321	CA	ALA	A	554	-2.711	75.834	38.289	1.00	21.10	6
4322	C	ALA	A	554	-1.169	75.788	38.321	1.00	17.79	6
4323	O	ALA	A	554	-0.631	75.032	39.141	1.00	17.24	8
4324	CB	ALA	A	554	-3.075	76.757	39.471	1.00	24.10	6
4325	N	ALA	A	555	-0.537	76.591	37.500	1.00	14.57	7
4326	CA	ALA	A	555	0.947	76.567	37.489	1.00	13.27	6
4327	C	ALA	A	555	1.568	77.356	38.600	1.00	13.78	6
4328	O	ALA	A	555	1.051	78.305	39.204	1.00	14.85	8
4329	CB	ALA	A	555	1.375	77.200	36.142	1.00	16.26	6
4330	N	GLY	A	556	2.874	76.966	38.857	1.00	12.02	7
4331	CA	GLY	A	556	3.602	76.639	39.947	1.00	14.17	6
4332	C	GLY	A	556	4.312	76.639	40.815	1.00	12.79	6
4333	O	GLY	A	556	4.121	75.415	40.670	1.00	12.92	8
4334	N	LEU	A	557	5.203	77.130	41.670	1.00	14.81	7
4335	CA	LEU	A	557	5.876	76.232	42.626	1.00	13.54	6
4336	C	LEU	A	557	4.918	75.806	43.737	1.00	12.41	6
4337	O	LEU	A	557	4.110	76.659	44.188	1.00	14.95	8
4338	CB	LEU	A	557	7.091	76.999	43.221	1.00	13.41	6
4339	CG	LEU	A	557	8.018	76.123	44.069	1.00	15.11	6
4340	CD1	LEU	A	557	8.847	75.220	43.141	1.00	13.33	6
4341	CD2	LEU	A	557	8.965	77.046	44.898	1.00	17.39	6
4342	N	THR	A	558	4.801	74.500	43.858	1.00	13.24	7
4343	CA	THR	A	558	3.807	73.941	44.818	1.00	13.26	6
4344	C	THR	A	558	4.469	72.927	45.741	1.00	16.19	6
4345	O	THR	A	558	5.520	72.402	45.379	1.00	16.82	8
4346	CB	THR	A	558	2.616	73.378	44.044	1.00	17.75	6
4347	OG1	THR	A	558	1.566	73.153	45.010	1.00	18.46	8
4348	CG2	THR	A	558	2.886	72.099	43.320	1.00	14.65	6
4349	N	ASP	A	559	3.750	72.559	46.821	1.00	13.53	7
4350	CA	ASP	A	559	4.351	71.569	47.739	1.00	13.26	6
4351	C	ASP	A	559	3.718	70.202	47.564	1.00	17.40	6
4352	O	ASP	A	559	2.469	70.015	47.579	1.00	17.26	8
4353	CB	ASP	A	559	4.129	72.027	49.195	1.00	13.77	6
4354	CG	ASP	A	559	4.998	73.219	49.580	1.00	25.94	6
4355	OD1	ASP	A	559	6.174	73.267	49.201	1.00	23.88	8
4356	OD2	ASP	A	559	4.468	74.127	50.251	1.00	28.55	8
4357	N	VAL	A	560	4.576	69.205	47.465	1.00	12.44	7
4358	CA	VAL	A	560	4.161	67.801	47.392	1.00	11.28	6
4359	C	VAL	A	560	4.193	67.207	48.826	1.00	14.98	6
4360	O	VAL	A	560	5.085	67.470	49.616	1.00	15.44	8
4361	CB	VAL	A	560	5.144	66.953	46.555	1.00	11.93	6
4362	CG1	VAL	A	560	4.738	65.496	46.488	1.00	16.58	6
4363	CG2	VAL	A	560	5.186	67.516	45.122	1.00	14.73	6
4364	N	LYS	A	561	3.136	66.439	49.097	1.00	15.49	7
4365	CA	LYS	A	561	2.989	65.835	50.443	1.00	14.90	6
4366	C	LYS	A	561	2.427	64.452	50.269	1.00	15.23	8
4367	O	LYS	A	561	1.502	64.157	49.522	1.00	16.90	6
4368	CB	LYS	A	561	1.960	66.720	51.206	1.00	18.50	6
4369	CG	LYS	A	561	1.847	66.177	52.656	1.00	22.42	6
4370	CD	LYS	A	561	1.025	67.110	53.523	1.00	25.69	6
4371	CE	LYS	A	561	-0.461	66.912	53.312	1.00	33.13	6
4372	NZ	LYS	A	561	-1.198	68.004	54.033	1.00	37.04	7
4373	N	VAL	A	562	2.947	63.489	51.047	1.00	16.07	7
4374	CA	VAL	A	562	2.525	62.108	51.051	1.00	14.01	6
4375	C	VAL	A	562	1.751	61.827	52.382	1.00	14.81	6
4376	O	VAL	A	562	2.150	62.324	53.398	1.00	17.73	8
4377	CB	VAL	A	562	3.735	61.158	51.036	1.00	16.60	6
4378	CG1	VAL	A	562	3.312	59.677	51.185	1.00	16.48	6
4379	CG2	VAL	A	562	4.473	61.228	49.672	1.00	17.80	8
4380	N	THR	A	563	0.603	61.156	52.140	1.00	17.67	7
4381	CA	THR	A	563	-0.181	60.806	53.380	1.00	18.18	6
4382	C	THR	A	563	-0.261	59.308	53.412	1.00	18.30	6
4383	O	THR	A	563	-0.679	58.642	52.456	1.00	19.93	8
4384	CB	THR	A	563	-1.542	61.496	53.361	1.00	18.11	6
4385	OG1	THR	A	563	-1.367	62.891	53.406	1.00	19.44	8

#	Atom	Res	Chain	ResNo	X	Y	Z	Occ	B	El
4386	CG2	THR	A	563	-2.281	61.112	54.705	1.00	20.36	6
4387	N	ALA	A	564	0.154	58.701	54.548	1.00	21.05	7
4388	CA	ALA	A	564	0.245	57.258	54.649	1.00	24.26	6
4389	C	ALA	A	564	-0.234	56.845	56.060	1.00	23.28	6
4390	O	ALA	A	564	0.104	57.536	57.008	1.00	21.58	8
4391	CB	ALA	A	564	1.658	56.711	54.437	1.00	25.03	6
4392	N	GLY	A	565	-1.218	55.968	56.140	1.00	30.08	7
4393	CA	GLY	A	565	-1.857	55.701	57.443	1.00	31.09	6
4394	C	GLY	A	565	-2.488	56.887	58.121	1.00	35.06	6
4395	O	GLY	A	565	-2.493	56.983	59.363	1.00	31.36	8
4396	N	GLY	A	566	-3.025	57.873	57.403	1.00	32.22	7
4397	CA	GLY	A	566	-3.549	59.083	58.011	1.00	31.70	6
4398	C	GLY	A	566	-2.523	60.107	58.438	1.00	31.07	6
4399	O	GLY	A	566	-2.936	61.213	58.829	1.00	32.69	8
4400	N	VAL	A	567	-1.202	59.887	58.271	1.00	26.17	7
4401	CA	VAL	A	567	-0.186	60.798	58.776	1.00	21.45	6
4402	C	VAL	A	567	0.537	61.488	57.557	1.00	17.95	6
4403	O	VAL	A	567	0.692	60.707	56.658	1.00	18.75	8
4404	CB	VAL	A	567	0.945	60.023	59.500	1.00	26.89	6
4405	CG1	VAL	A	567	1.981	60.946	60.096	1.00	30.00	6
4406	CG2	VAL	A	567	0.308	59.088	60.567	1.00	33.86	6
4407	N	SER	A	568	0.687	62.774	57.699	1.00	19.77	7
4408	CA	SER	A	568	1.267	63.410	56.486	1.00	19.16	6
4409	C	SER	A	568	2.770	63.585	56.688	1.00	19.28	6
4410	O	SER	A	568	3.351	63.752	57.763	1.00	21.35	8
4411	CB	SER	A	568	0.574	64.705	56.156	1.00	29.69	6
4412	OG	SER	A	568	0.595	65.549	57.266	1.00	41.18	8
4413	N	SER	A	569	3.399	63.581	55.503	1.00	19.22	7
4414	CA	SER	A	569	4.867	63.784	55.480	1.00	17.43	6
4415	C	SER	A	569	5.229	65.245	55.568	1.00	17.17	6
4416	O	SER	A	569	4.519	66.266	55.502	1.00	17.53	8
4417	CB	SER	A	569	5.381	63.220	54.137	1.00	17.58	6
4418	OG	SER	A	569	5.066	64.083	53.025	1.00	15.50	8
4419	N	ASN	A	570	6.572	65.473	55.538	1.00	14.77	7
4420	CA	ASN	A	570	7.143	66.776	55.253	1.00	13.84	6
4421	C	ASN	A	570	6.848	67.107	53.752	1.00	13.33	6
4422	O	ASN	A	570	6.527	66.217	52.996	1.00	16.50	8
4423	CB	ASN	A	570	8.670	66.723	55.435	1.00	17.22	6
4424	CG	ASN	A	570	9.363	65.581	54.758	1.00	16.95	6
4425	OD1	ASN	A	570	9.038	64.399	54.745	1.00	15.12	8
4426	ND2	ASN	A	570	10.455	65.853	54.023	1.00	16.51	7
4427	N	LEU	A	571	7.108	68.381	53.486	1.00	14.82	7
4428	CA	LEU	A	571	6.705	68.878	52.131	1.00	12.49	6
4429	C	LEU	A	571	7.959	69.019	51.282	1.00	15.87	6
4430	O	LEU	A	571	9.024	69.355	51.828	1.00	16.79	8
4431	CB	LEU	A	571	6.147	70.286	52.340	1.00	15.18	6
4432	CG	LEU	A	571	4.911	70.396	53.250	1.00	22.77	6
4433	CD1	LEU	A	571	4.368	71.807	53.300	1.00	22.26	6
4434	CD2	LEU	A	571	3.834	69.448	52.757	1.00	22.04	6
4435	N	TYR	A	572	7.747	68.878	49.946	1.00	13.65	7
4436	CA	TYR	A	572	8.930	69.100	49.067	1.00	12.87	6
4437	C	TYR	A	572	8.322	69.878	47.880	1.00	14.52	6
4438	O	TYR	A	572	7.347	69.399	47.262	1.00	14.71	8
4439	CB	TYR	A	572	9.480	67.704	48.701	1.00	12.62	6
4440	CG	TYR	A	572	10.887	67.679	48.121	1.00	13.61	6
4441	CD1	TYR	A	572	11.079	68.175	46.835	1.00	13.57	6
4442	CD2	TYR	A	572	11.946	67.161	48.845	1.00	15.40	6
4443	CE1	TYR	A	572	12.361	68.168	46.257	1.00	13.80	6
4444	CE2	TYR	A	572	13.221	67.135	48.283	1.00	12.70	6
4445	CZ	TYR	A	572	13.400	67.629	47.002	1.00	14.33	6
4446	OH	TYR	A	572	14.710	67.607	46.466	1.00	13.40	8
4447	N	SER	A	573	9.060	70.874	47.387	1.00	13.37	7
4448	CA	SER	A	573	8.571	71.734	46.294	1.00	12.21	6
4449	C	SER	A	573	8.665	71.055	44.920	1.00	12.75	6
4450	O	SER	A	573	9.520	70.234	44.703	1.00	12.64	8
4451	CB	ASER	A	573	9.436	72.999	46.162	0.60	17.33	6
4452	OG	ASER	A	573	9.459	73.767	47.331	0.60	22.77	8
4452	CB	BSER	A	573	9.408	73.025	46.276	0.40	15.40	6
4453	OG	BSER	A	573	10.793	72.728	46.149	0.40	16.17	8
4453	N	TYR	A	574	7.838	71.568	44.000	1.00	11.56	7
4454	CA	TYR	A	574	7.912	71.077	42.604	1.00	10.89	6
4455	C	TYR	A	574	7.374	72.249	41.771	1.00	13.28	6
4456	O	TYR	A	574	6.344	72.860	42.138	1.00	12.17	8
4457	CB	TYR	A	574	7.041	69.857	42.405	1.00	11.94	6
4458	CG	TYR	A	574	6.917	69.379	40.971	1.00	11.37	6
4459	CD1	TYR	A	574	7.921	68.521	40.467	1.00	12.69	6
4460	CD2	TYR	A	574	5.867	69.781	40.162	1.00	11.85	6
4461	CE1	TYR	A	574	7.863	68.052	39.154	1.00	13.23	6
4462	CE2	TYR	A	574	5.792	69.313	38.366	1.00	13.93	6
4463	CZ	TYR	A	574	6.796	68.466	38.366	1.00	11.32	6
4464	OH	TYR	A	574	6.679	68.006	37.078	1.00	12.54	8
4465	N	ASN	A	575	7.992	72.510	40.627	1.00	12.18	7
4466	CA	ASN	A	575	7.578	73.682	39.820	1.00	12.18	6
4467	C	ASN	A	575	6.738	73.306	38.608	1.00	12.59	6

Serial	Atom	Res	Ch	Seq	X	Y	Z	Occ	B	El
4468	O	ASN	A	575	7.171	72.699	37.598	1.00	12.20	8
4469	CB	ASN	A	575	8.898	74.351	39.331	1.00	12.02	6
4470	CG	ASN	A	575	8.635	75.707	38.700	1.00	17.00	6
4471	OD1	ASN	A	575	7.562	76.292	38.872	1.00	14.54	8
4472	ND2	ASN	A	575	9.608	76.223	37.934	1.00	13.02	7
4473	N	ILE	A	576	5.400	73.515	38.739	1.00	10.95	7
4474	CA	ILE	A	576	4.506	73.221	37.604	1.00	10.36	6
4475	C	ILE	A	576	4.485	74.374	36.602	1.00	12.05	6
4476	O	ILE	A	576	4.145	75.497	36.930	1.00	13.39	8
4477	CB	ILE	A	576	3.036	71.956	38.127	1.00	12.95	6
4478	CG1	ILE	A	576	3.082	71.956	39.190	1.00	13.10	6
4479	CG2	ILE	A	576	2.079	72.677	36.966	1.00	14.18	6
4480	CD1	ILE	A	576	1.709	71.488	39.691	1.00	15.07	6
4481	N	LEU	A	577	4.883	74.054	35.368	1.00	10.51	7
4482	CA	LEU	A	577	4.908	75.052	34.303	1.00	12.74	6
4483	C	LEU	A	577	3.480	75.308	33.772	1.00	11.80	6
4484	O	LEU	A	577	2.572	74.474	34.021	1.00	13.37	8
4485	CB	LEU	A	577	5.757	74.463	33.137	1.00	11.48	6
4486	CG	LEU	A	577	7.226	74.238	33.569	1.00	12.21	6
4487	CD1	LEU	A	577	7.982	73.601	32.421	1.00	13.03	6
4488	CD2	LEU	A	577	7.897	75.576	33.989	1.00	13.85	6
4489	N	SER	A	578	3.398	76.385	33.006	1.00	11.76	7
4490	CA	SER	A	578	2.037	76.726	32.475	1.00	15.00	6
4491	C	SER	A	578	1.681	75.960	31.215	1.00	17.37	6
4492	O	SER	A	578	0.553	76.134	30.685	1.00	17.93	8
4493	CB	SER	A	578	2.081	78.230	32.132	1.00	13.25	6
4494	OG	SER	A	578	2.321	78.967	33.305	1.00	14.45	8
4495	N	GLY	A	579	2.539	75.095	30.688	1.00	15.28	7
4496	CA	GLY	A	579	2.266	74.270	29.497	1.00	14.88	6
4497	C	GLY	A	579	3.627	74.010	28.818	1.00	13.39	6
4498	O	GLY	A	579	4.670	74.450	29.318	1.00	14.78	8
4499	N	THR	A	580	3.518	73.227	27.740	1.00	11.35	7
4500	CA	THR	A	580	4.803	72.961	27.005	1.00	13.16	6
4501	C	THR	A	580	5.419	74.292	26.647	1.00	14.61	6
4502	O	THR	A	580	4.747	75.282	26.349	1.00	14.26	8
4503	CB	THR	A	580	4.517	71.676	25.835	1.00	13.57	6
4504	OG1	THR	A	580	5.753	72.715	25.176	1.00	16.06	8
4505	CG2	THR	A	580	3.688	72.715	24.732	1.00	20.75	6
4506	N	GLN	A	581	6.787	74.346	26.706	1.00	11.89	7
4507	CA	GLN	A	581	7.454	75.658	26.609	1.00	10.06	6
4508	C	GLN	A	581	8.012	75.905	25.148	1.00	10.19	6
4509	O	GLN	A	581	8.163	75.020	24.338	1.00	12.94	8
4510	CB	GLN	A	581	8.751	75.601	27.465	1.00	11.27	6
4511	CG	GLN	A	581	8.407	75.301	28.975	1.00	10.49	6
4512	CD	GLN	A	581	7.920	76.579	29.654	1.00	12.45	6
4513	OE1	GLN	A	581	8.545	77.588	29.921	1.00	13.44	8
4514	NE2	GLN	A	581	6.582	76.528	30.002	1.00	11.11	7
4515	N	THR	A	582	8.202	77.209	24.953	1.00	9.52	7
4516	CA	THR	A	582	8.978	77.646	23.772	1.00	10.13	6
4517	C	THR	A	582	10.049	78.609	24.293	1.00	10.70	6
4518	O	THR	A	582	9.921	79.235	25.313	1.00	12.73	8
4519	CB	THR	A	582	8.018	78.306	22.763	1.00	12.88	6
4520	OG1	THR	A	582	8.736	78.798	21.599	1.00	12.92	8
4521	CG2	THR	A	582	7.265	79.504	23.359	1.00	11.55	6
4522	N	SER	A	583	11.166	78.712	23.507	1.00	11.67	7
4523	CA	SER	A	583	12.321	79.527	23.931	1.00	12.58	6
4524	C	SER	A	583	12.300	80.783	23.061	1.00	11.58	6
4525	O	SER	A	583	12.496	80.687	21.844	1.00	12.44	8
4526	CB	SER	A	583	13.612	78.697	23.747	1.00	12.53	6
4527	OG	SER	A	583	14.755	79.449	24.240	1.00	14.73	8
4528	N	VAL	A	584	12.118	81.899	23.781	1.00	10.19	7
4529	CA	VAL	A	584	11.741	83.141	23.016	1.00	9.76	6
4530	C	VAL	A	584	12.721	84.268	23.298	1.00	11.27	6
4531	O	VAL	A	584	13.066	84.570	24.440	1.00	11.87	8
4532	CB	VAL	A	584	10.358	83.645	23.506	1.00	12.33	6
4533	CG1	VAL	A	584	10.041	84.968	22.757	1.00	14.53	6
4534	CG2	VAL	A	584	9.279	82.604	23.220	1.00	12.71	6
4535	N	VAL	A	585	13.307	84.868	22.244	1.00	10.77	7
4536	CA	VAL	A	585	14.166	86.030	22.421	1.00	9.69	6
4537	C	VAL	A	585	13.244	87.219	22.693	1.00	12.06	6
4538	O	VAL	A	585	12.450	87.576	21.803	1.00	14.60	8
4539	CB	VAL	A	585	14.882	86.310	21.064	1.00	11.53	6
4540	CG1	VAL	A	585	15.712	87.605	21.158	1.00	15.03	6
4541	CG2	VAL	A	585	15.750	85.111	20.717	1.00	14.67	6
4542	N	PHE	A	586	13.361	87.762	23.908	1.00	12.32	7
4543	CA	PHE	A	586	12.622	88.986	24.243	1.00	12.19	6
4544	C	PHE	A	586	13.584	90.161	24.063	1.00	14.77	6
4545	O	PHE	A	586	14.668	90.209	24.709	1.00	12.38	8
4546	CB	PHE	A	586	12.174	88.944	25.704	1.00	12.57	6
4547	CG	PHE	A	586	10.886	88.184	25.987	1.00	11.39	6
4548	CD1	PHE	A	586	10.879	86.809	26.009	1.00	13.56	6
4549	CD2	PHE	A	586	9.694	88.868	26.231	1.00	13.70	6
4550	CE1	PHE	A	586	9.758	86.060	26.284	1.00	13.86	6
4551	CE2	PHE	A	586	8.543	88.102	26.520	1.00	12.32	6

4552	CZ	PHE	A	586	8.577	86.707	26.558	1.00	11.70	6
4553	N	THR	A	587	13.189	91.146	23.239	1.00	14.16	7
4554	CA	THR	A	587	14.016	92.324	23.066	1.00	12.96	6
4555	C	THR	A	587	13.247	93.561	23.468	1.00	13.58	6
4556	O	THR	A	587	12.072	93.651	23.058	1.00	15.22	8
4557	CB	THR	A	587	14.421	92.464	21.555	1.00	11.66	6
4558	OG1	THR	A	587	15.145	91.299	21.141	1.00	15.06	8
4559	CG2	THR	A	587	15.331	93.708	21.318	1.00	13.37	6
4560	N	VAL	A	588	13.829	94.487	24.195	1.00	14.03	7
4561	CA	VAL	A	588	13.156	95.779	24.469	1.00	14.79	6
4562	C	VAL	A	588	14.079	96.867	23.912	1.00	14.40	6
4563	O	VAL	A	588	15.258	97.015	24.225	1.00	15.56	8
4564	CB	VAL	A	588	12.863	95.933	25.971	1.00	14.24	6
4565	CG1	VAL	A	588	14.111	95.935	26.870	1.00	14.60	6
4566	CG2	VAL	A	588	12.079	97.255	26.232	1.00	13.38	6
4567	N	LYS	A	589	13.478	97.570	22.915	1.00	15.39	7
4568	CA	LYS	A	589	14.212	98.670	22.258	1.00	16.20	6
4569	C	LYS	A	589	14.180	100.009	22.953	1.00	18.01	6
4570	O	LYS	A	589	13.230	100.355	23.652	1.00	15.39	8
4571	CB	LYS	A	589	13.597	98.820	20.860	1.00	16.23	6
4572	CG	LYS	A	589	13.908	97.588	20.008	1.00	17.01	6
4573	CD	LYS	A	589	13.275	97.825	18.634	1.00	24.71	6
4574	CE	LYS	A	589	13.494	96.582	17.792	1.00	37.49	6
4575	NZ	LYS	A	589	13.368	96.852	16.321	1.00	51.36	7
4576	N	SER	A	590	15.302	100.747	22.828	1.00	16.34	7
4577	CA	SER	A	590	15.371	102.116	23.293	1.00	19.32	6
4578	C	SER	A	590	15.006	102.295	24.746	1.00	19.68	6
4579	O	SER	A	590	14.185	103.146	25.151	1.00	17.24	8
4580	CB	SER	A	590	15.006	103.009	22.421	1.00	20.32	6
4581	OG	SER	A	590	14.867	102.933	21.046	1.00	22.81	8
4582	N	ALA	A	591	15.698	101.514	25.612	1.00	17.08	7
4583	CA	ALA	A	591	15.458	101.617	27.048	1.00	16.02	6
4584	C	ALA	A	591	16.178	102.871	27.530	1.00	18.06	6
4585	O	ALA	A	591	17.152	103.325	26.878	1.00	15.85	8
4586	CB	ALA	A	591	16.045	100.326	27.695	1.00	15.17	6
4587	N	PRO	A	592	15.872	103.303	28.730	1.00	18.28	7
4588	CA	PRO	A	592	16.493	104.529	29.298	1.00	18.18	6
4589	C	PRO	A	592	17.967	104.319	29.512	1.00	22.69	6
4590	O	PRO	A	592	18.463	103.180	29.693	1.00	19.12	8
4591	CB	PRO	A	592	15.762	104.797	30.621	1.00	19.23	6
4592	CG	PRO	A	592	14.433	104.085	30.387	1.00	23.18	6
4593	CD	PRO	A	592	14.795	102.833	29.585	1.00	17.71	6
4594	N	PRO	A	593	18.778	105.367	29.550	1.00	21.02	7
4595	CA	PRO	A	593	20.204	105.252	29.779	1.00	20.83	6
4596	C	PRO	A	593	20.494	104.537	31.094	1.00	20.17	6
4597	O	PRO	A	593	19.811	104.784	32.102	1.00	21.50	8
4598	CB	PRO	A	593	20.715	106.727	29.941	1.00	22.70	6
4599	CG	PRO	A	593	19.643	107.491	29.165	1.00	24.71	6
4600	CD	PRO	A	593	18.337	106.774	29.432	1.00	23.44	6
4601	N	THR	A	594	21.530	103.723	31.165	1.00	19.91	7
4602	CA	THR	A	594	21.909	103.060	32.402	1.00	21.78	6
4603	C	THR	A	594	23.380	103.366	32.686	1.00	23.13	6
4604	O	THR	A	594	24.138	103.737	31.787	1.00	23.96	8
4605	CB	THR	A	594	21.729	101.521	32.292	1.00	23.27	6
4606	OG1	THR	A	594	22.466	101.094	31.140	1.00	20.06	8
4607	CG2	THR	A	594	20.245	101.183	32.157	1.00	21.33	6
4608	N	ASN	A	595	20.764	103.210	33.928	1.00	23.96	7
4609	CA	ASN	A	595	25.142	103.193	34.370	1.00	28.93	6
4610	C	ASN	A	595	25.614	101.791	34.716	1.00	31.88	6
4611	O	ASN	A	595	24.847	100.849	34.915	1.00	22.74	8
4612	CB	ASN	A	595	25.285	104.099	35.620	1.00	34.82	6
4613	CG	ASN	A	595	24.970	105.532	35.180	1.00	36.76	6
4614	OD1	ASN	A	595	24.047	106.190	35.647	1.00	42.57	8
4615	ND2	ASN	A	595	25.738	106.004	34.208	1.00	38.16	7
4616	N	LEU	A	596	26.939	101.654	34.839	1.00	34.83	7
4617	CA	LEU	A	596	27.559	100.380	35.212	1.00	37.55	6
4618	C	LEU	A	596	26.947	99.803	36.488	1.00	31.76	6
4619	O	LEU	A	596	26.589	100.502	37.435	1.00	35.49	8
4620	CB	LEU	A	596	29.051	100.642	35.407	1.00	48.79	6
4621	CG	LEU	A	596	30.042	99.528	35.681	1.00	52.60	6
4622	CD1	LEU	A	596	29.894	98.971	37.092	1.00	56.16	6
4623	CD2	LEU	A	596	29.934	98.433	34.628	1.00	56.63	6
4624	N	GLY	A	597	26.492	98.556	36.348	1.00	30.34	7
4625	CA	GLY	A	597	25.861	97.860	37.473	1.00	30.50	6
4626	C	GLY	A	597	24.337	97.822	37.288	1.00	25.20	6
4627	O	GLY	A	597	23.705	96.996	37.949	1.00	23.57	8
4628	N	ASP	A	598	23.780	98.803	36.544	1.00	21.92	7
4629	CA	ASP	A	598	22.315	98.799	36.399	1.00	18.64	6
4630	C	ASP	A	598	21.982	97.666	35.429	1.00	21.53	6
4631	O	ASP	A	598	22.702	97.563	34.399	1.00	20.69	8
4632	CB	ASP	A	598	21.814	100.096	35.763	1.00	17.61	6
4633	CG	ASP	A	598	22.046	101.382	36.536	1.00	20.17	6
4634	OD1	ASP	A	598	22.364	101.347	37.751	1.00	19.25	8
4635	OD2	ASP	A	598	21.858	102.432	35.868	1.00	22.20	8

4636	N	LYS	A	599	20.861	96.919	35.612	1.00	18.66	7
4637	CA	LYS	A	599	20.593	95.891	34.605	1.00	15.76	6
4638	C	LYS	A	599	19.056	95.824	34.461	1.00	11.90	6
4639	O	LYS	A	599	18.320	96.237	35.358	1.00	19.04	8
4640	CB	LYS	A	599	21.049	94.484	35.006	1.00	23.86	6
4641	CG	LYS	A	599	22.597	94.419	35.031	1.00	25.32	6
4642	CD	LYS	A	599	23.118	93.007	35.252	1.00	26.43	6
4643	CE	LYS	A	599	24.656	93.063	35.120	1.00	32.43	6
4644	NZ	LYS	A	599	25.179	91.656	35.106	1.00	35.89	7
4645	N	ILE	A	600	18.623	95.442	33.287	1.00	13.51	7
4646	CA	ILE	A	600	17.178	95.336	33.021	1.00	12.20	6
4647	C	ILE	A	600	16.746	93.886	33.130	1.00	13.74	6
4648	O	ILE	A	600	17.476	92.927	32.799	1.00	14.24	8
4649	CB	ILE	A	600	16.938	95.829	31.554	1.00	13.87	6
4650	CG1	ILE	A	600	17.249	97.335	31.566	1.00	21.52	6
4651	CG2	ILE	A	600	15.499	95.618	31.051	1.00	15.85	6
4652	CD1	ILE	A	600	16.798	98.104	30.313	1.00	23.52	6
4653	N	TYR	A	601	15.545	93.745	33.676	1.00	13.09	7
4654	CA	TYR	A	601	14.926	92.453	33.883	1.00	12.76	6
4655	C	TYR	A	601	13.531	92.475	33.288	1.00	13.27	6
4656	O	TYR	A	601	12.914	93.491	32.986	1.00	15.09	8
4657	CB	TYR	A	601	14.751	92.119	35.387	1.00	14.11	6
4658	CG	TYR	A	601	16.065	91.805	36.078	1.00	11.98	6
4659	CD1	TYR	A	601	16.902	92.834	36.502	1.00	12.52	6
4660	CD2	TYR	A	601	16.473	90.490	36.242	1.00	12.95	6
4661	CE1	TYR	A	601	18.138	92.544	37.097	1.00	13.54	6
4662	CE2	TYR	A	601	17.673	90.213	36.888	1.00	14.26	6
4663	CZ	TYR	A	601	18.499	91.239	37.276	1.00	15.84	6
4664	OH	TYR	A	601	19.691	90.903	37.935	1.00	16.92	8
4665	N	LEU	A	602	12.986	91.285	33.068	1.00	11.99	7
4666	CA	LEU	A	602	11.657	91.057	32.494	1.00	13.20	6
4667	C	LEU	A	602	10.762	90.434	33.580	1.00	14.26	6
4668	O	LEU	A	602	11.148	89.522	34.284	1.00	13.06	8
4669	CB	LEU	A	602	11.794	90.060	31.334	1.00	13.29	6
4670	CG	LEU	A	602	10.506	89.432	30.800	1.00	10.59	6
4671	CD1	LEU	A	602	9.579	90.481	30.155	1.00	14.08	6
4672	CD2	LEU	A	602	10.864	88.385	29.704	1.00	11.83	6
4673	N	THR	A	603	9.482	90.861	33.649	1.00	12.02	7
4674	CA	THR	A	603	8.570	90.233	34.635	1.00	12.81	6
4675	C	THR	A	603	7.158	90.267	34.066	1.00	12.50	6
4676	O	THR	A	603	6.876	91.096	33.183	1.00	14.59	8
4677	CB	THR	A	603	8.711	90.983	35.969	1.00	14.38	6
4678	OG1	THR	A	603	7.971	90.241	36.987	1.00	14.53	8
4679	CG2	THR	A	603	8.152	92.414	35.971	1.00	15.35	6
4680	N	GLY	A	604	6.310	89.331	34.489	1.00	11.51	7
4681	CA	GLY	A	604	4.995	89.274	33.787	1.00	11.82	6
4682	C	GLY	A	604	4.042	88.348	34.530	1.00	12.03	6
4683	O	GLY	A	604	4.358	87.909	35.636	1.00	13.68	8
4684	N	ASN	A	605	2.911	88.082	33.859	1.00	12.03	7
4685	CA	ASN	A	605	1.782	87.516	34.637	1.00	11.85	6
4686	C	ASN	A	605	1.680	86.017	34.575	1.00	12.50	6
4687	O	ASN	A	605	0.615	85.441	34.625	1.00	14.07	8
4688	CB	ASN	A	605	0.481	88.161	34.031	1.00	13.77	6
4689	CG	ASN	A	605	0.265	87.610	32.629	1.00	17.19	6
4690	OD1	ASN	A	605	1.080	87.101	31.828	1.00	13.63	8
4691	ND2	ASN	A	605	-1.025	87.681	32.165	1.00	16.36	7
4692	N	ILE	A	606	2.823	85.308	34.560	1.00	13.26	7
4693	CA	ILE	A	606	2.836	83.860	34.628	1.00	11.73	6
4694	C	ILE	A	606	4.099	83.493	35.448	1.00	13.02	6
4695	O	ILE	A	606	5.005	84.349	35.511	1.00	13.34	8
4696	CB	ILE	A	606	2.995	83.184	33.259	1.00	11.81	6
4697	CG1	ILE	A	606	4.029	83.874	32.346	1.00	12.94	6
4698	CG2	ILE	A	606	1.625	83.138	32.550	1.00	15.56	6
4699	CD1	ILE	A	606	4.300	82.981	31.113	1.00	15.13	6
4700	N	PRO	A	607	4.181	82.331	36.033	1.00	13.09	7
4701	CA	PRO	A	607	5.294	81.982	36.937	1.00	12.68	6
4702	C	PRO	A	607	6.616	81.880	36.189	1.00	10.98	6
4703	O	PRO	A	607	7.700	82.258	36.671	1.00	14.73	8
4704	CB	PRO	A	607	4.895	80.647	37.607	1.00	13.69	6
4705	CG	PRO	A	607	3.818	80.118	36.675	1.00	16.48	6
4706	CD	PRO	A	607	3.081	81.335	36.098	1.00	16.43	6
4707	N	GLU	A	608	6.531	81.525	34.880	1.00	11.74	7
4708	CA	GLU	A	608	7.747	81.559	34.031	1.00	10.58	6
4709	C	GLU	A	608	8.382	82.958	33.994	1.00	13.60	6
4710	O	GLU	A	608	9.606	83.060	33.731	1.00	11.79	8
4711	CB	GLU	A	608	7.437	81.056	32.616	1.00	11.09	6
4712	CG	GLU	A	608	7.145	79.550	32.522	1.00	10.94	6
4713	CD	GLU	A	608	5.674	79.163	32.699	1.00	14.31	6
4714	OE1	GLU	A	608	4.875	79.973	33.221	1.00	12.87	8
4715	OE2	GLU	A	608	5.344	78.040	32.280	1.00	13.22	8
4716	N	LEU	A	609	7.571	83.999	34.112	1.00	11.31	7
4717	CA	LEU	A	609	8.093	85.359	34.095	1.00	11.00	6
4718	C	LEU	A	609	7.944	86.002	35.453	1.00	12.01	6
4719	O	LEU	A	609	7.959	87.224	35.602	1.00	14.24	8

100

No	Atom	Res	Ch	Num	X	Y	Z	Occ	B	E
4720	CB	LEU	A	609	7.249	86.188	33.078	1.00	14.30	6
4721	CG	LEU	A	609	7.424	85.668	31.627	1.00	13.57	6
4722	CD1	LEU	A	609	6.672	86.643	30.695	1.00	18.28	6
4723	CD2	LEU	A	609	8.912	85.648	31.238	1.00	14.71	6
4724	N	GLY	A	610	7.854	85.194	36.525	1.00	10.78	7
4725	CA	GLY	A	610	7.950	85.729	37.864	1.00	10.94	6
4726	C	GLY	A	610	6.664	86.230	38.554	1.00	13.37	6
4727	O	GLY	A	610	6.767	86.832	39.647	1.00	13.90	8
4728	N	ASN	A	611	5.525	86.174	37.906	1.00	13.56	7
4729	CA	ASN	A	611	4.262	86.696	38.463	1.00	14.13	6
4730	C	ASN	A	611	4.462	88.079	39.072	1.00	13.68	6
4731	O	ASN	A	611	4.063	88.427	40.207	1.00	14.78	8
4732	CB	ASN	A	611	3.695	85.782	39.555	1.00	13.73	6
4733	CG	ASN	A	611	3.110	84.538	38.890	1.00	16.98	6
4734	OD1	ASN	A	611	2.353	84.585	37.897	1.00	17.26	8
4735	ND2	ASN	A	611	3.501	83.420	39.478	1.00	16.08	7
4736	N	TRP	A	612	5.010	88.962	38.242	1.00	13.00	7
4737	CA	TRP	A	612	5.298	90.339	38.534	1.00	13.35	6
4738	C	TRP	A	612	6.420	90.676	39.608	1.00	14.25	6
4739	O	TRP	A	612	6.420	91.775	40.107	1.00	15.26	8
4740	CB	TRP	A	612	3.915	91.044	38.832	1.00	13.65	6
4741	CG	TRP	A	612	2.965	91.036	37.664	1.00	13.25	6
4742	CD1	TRP	A	612	1.775	90.359	37.632	1.00	13.84	6
4743	CD2	TRP	A	612	3.138	91.564	36.359	1.00	14.07	6
4744	NE1	TRP	A	612	1.160	90.489	36.402	1.00	13.90	7
4745	CE2	TRP	A	612	1.983	91.288	35.617	1.00	16.64	6
4746	CE3	TRP	A	612	4.148	92.402	35.815	1.00	15.91	6
4747	CZ2	TRP	A	612	1.843	91.676	34.288	1.00	14.35	6
4748	CZ3	TRP	A	612	3.980	92.856	34.506	1.00	14.55	6
4749	CH2	TRP	A	612	2.844	92.490	33.747	1.00	14.05	6
4750	N	SER	A	613	7.114	89.674	40.049	1.00	14.59	7
4751	CA	SER	A	613	8.136	89.943	41.025	1.00	16.76	6
4752	C	SER	A	613	9.171	90.916	40.471	1.00	15.94	6
4753	O	SER	A	613	9.510	90.841	39.281	1.00	15.44	8
4754	CB	SER	A	613	8.804	88.572	41.315	1.00	14.15	6
4755	OG	SER	A	613	9.842	88.719	42.250	1.00	15.35	8
4756	N	THR	A	614	9.821	91.648	41.406	1.00	13.40	7
4757	CA	THR	A	614	11.021	92.394	41.037	1.00	13.18	6
4758	C	THR	A	614	12.238	91.869	41.796	1.00	12.85	6
4759	O	THR	A	614	13.311	92.464	41.777	1.00	15.88	8
4760	CB	THR	A	614	10.895	93.904	41.297	1.00	16.41	6
4761	OG1	THR	A	614	10.626	94.073	42.724	1.00	18.07	8
4762	CG2	THR	A	614	9.718	94.523	40.520	1.00	18.09	6
4763	N	ASP	A	615	12.033	90.680	42.425	1.00	13.65	7
4764	CA	ASP	A	615	13.138	90.024	43.143	1.00	15.24	6
4765	C	ASP	A	615	14.084	89.416	42.093	1.00	12.74	6
4766	O	ASP	A	615	13.582	88.794	41.143	1.00	13.71	8
4767	CB	ASP	A	615	12.511	88.981	44.075	1.00	14.77	6
4768	CG	ASP	A	615	13.634	88.423	44.970	1.00	21.15	6
4769	OD1	ASP	A	615	13.956	88.932	46.048	1.00	23.65	8
4770	OD2	ASP	A	615	14.204	87.431	44.586	1.00	18.42	8
4771	N	THR	A	616	15.390	89.536	42.350	1.00	12.77	7
4772	CA	THR	A	616	16.387	89.067	41.365	1.00	13.57	6
4773	C	THR	A	616	17.391	88.146	42.072	1.00	15.88	6
4774	O	THR	A	616	18.487	87.902	41.566	1.00	14.89	8
4775	CB	THR	A	616	17.144	90.224	40.674	1.00	14.32	6
4776	OG1	THR	A	616	17.752	91.045	41.713	1.00	15.99	8
4777	CG2	THR	A	616	16.189	91.128	39.884	1.00	16.33	6
4778	N	SER	A	617	16.981	87.590	43.216	1.00	13.15	7
4779	CA	SER	A	617	17.813	86.641	43.956	1.00	12.49	6
4780	C	SER	A	617	17.759	85.265	43.312	1.00	13.68	6
4781	O	SER	A	617	17.193	85.034	42.235	1.00	12.79	8
4782	CB	SER	A	617	17.271	86.583	45.396	1.00	14.38	6
4783	OG	SER	A	617	16.056	85.882	45.441	1.00	15.10	8
4784	N	GLY	A	618	18.327	84.268	44.080	1.00	15.05	7
4785	CA	GLY	A	618	18.300	82.869	43.621	1.00	13.01	6
4786	C	GLY	A	618	16.967	82.147	43.811	1.00	13.49	6
4787	O	GLY	A	618	16.835	80.938	43.552	1.00	15.47	8
4788	N	ALA	A	619	15.893	82.847	44.200	1.00	13.99	7
4789	CA	ALA	A	619	14.592	82.218	44.333	1.00	14.93	6
4790	C	ALA	A	619	14.033	81.691	43.015	1.00	15.28	6
4791	O	ALA	A	619	14.489	82.043	41.907	1.00	15.28	8
4792	CB	ALA	A	619	13.582	83.263	44.854	1.00	17.52	6
4793	N	VAL	A	620	13.043	80.805	43.172	1.00	11.74	7
4794	CA	VAL	A	620	12.419	80.230	41.967	1.00	12.25	6
4795	C	VAL	A	620	11.259	81.126	41.502	1.00	12.14	6
4796	O	VAL	A	620	10.533	81.679	42.344	1.00	14.38	8
4797	CB	VAL	A	620	11.796	78.892	42.436	1.00	15.46	6
4798	CG1	VAL	A	620	11.242	78.231	41.173	1.00	13.86	6
4799	CG2	VAL	A	620	12.924	77.941	42.923	1.00	15.04	6
4800	N	ASN	A	621	11.157	81.254	40.168	1.00	12.23	7
4801	CA	ASN	A	621	10.066	82.044	39.571	1.00	12.45	6
4802	C	ASN	A	621	9.968	83.484	40.032	1.00	12.62	6
4803	O	ASN	A	621	8.860	83.896	40.480	1.00	12.87	8

4804	CB	ASN	A	621	8.676	81.346	39.732	1.00	10.99	6
4805	CG	ASN	A	621	8.656	79.972	39.084	1.00	11.13	6
4806	OD1	ASN	A	621	9.398	79.689	38.141	1.00	13.90	8
4807	ND2	ASN	A	621	7.742	79.105	39.596	1.00	13.32	7
4808	N	ASN	A	622	11.106	84.199	40.048	1.00	13.97	7
4809	CA	ASN	A	622	11.116	85.623	40.322	1.00	11.35	6
4810	C	ASN	A	622	11.418	86.389	39.037	1.00	12.47	6
4811	O	ASN	A	622	11.090	85.862	37.960	1.00	12.66	8
4812	CB	ASN	A	622	12.073	85.956	41.471	1.00	12.54	6
4813	CG	ASN	A	622	13.543	85.545	41.151	1.00	12.82	6
4814	OD1	ASN	A	622	13.811	85.093	40.044	1.00	12.70	8
4815	ND2	ASN	A	622	14.377	85.753	42.196	1.00	13.72	7
4816	N	ALA	A	623	11.883	87.626	39.098	1.00	11.27	7
4817	CA	ALA	A	623	12.088	88.339	37.812	1.00	12.38	6
4818	C	ALA	A	623	13.107	87.592	36.947	1.00	13.25	6
4819	O	ALA	A	623	14.035	87.018	37.469	1.00	13.43	8
4820	CB	ALA	A	623	12.586	89.748	38.120	1.00	14.46	6
4821	N	GLN	A	624	12.905	87.740	35.631	1.00	12.11	7
4822	CA	GLN	A	624	13.742	87.034	34.660	1.00	11.87	6
4823	C	GLN	A	624	14.828	87.912	34.071	1.00	12.70	6
4824	O	GLN	A	624	14.646	89.092	33.762	1.00	13.14	8
4825	CB	GLN	A	624	12.880	86.477	33.521	1.00	14.55	6
4826	CG	GLN	A	624	11.779	85.492	34.084	1.00	13.33	6
4827	CD	GLN	A	624	12.451	84.323	34.754	1.00	14.35	6
4828	OE1	GLN	A	624	13.213	83.546	34.096	1.00	15.81	8
4829	NE2	GLN	A	624	12.268	84.078	36.040	1.00	14.08	7
4830	N	GLY	A	625	15.989	87.264	33.941	1.00	12.93	7
4831	CA	GLY	A	625	17.183	87.950	33.418	1.00	12.61	6
4832	C	GLY	A	625	18.392	87.917	34.310	1.00	12.24	6
4833	O	GLY	A	625	18.497	86.888	34.934	1.00	13.82	8
4834	N	PRO	A	626	19.296	88.815	34.131	1.00	12.81	7
4835	CA	PRO	A	626	19.237	90.097	33.510	1.00	15.64	6
4836	C	PRO	A	626	19.288	89.998	32.003	1.00	15.42	6
4837	O	PRO	A	626	19.675	88.992	31.364	1.00	15.65	8
4838	CB	PRO	A	626	20.374	91.012	34.019	1.00	15.71	6
4839	CG	PRO	A	626	21.401	89.927	34.320	1.00	14.63	6
4840	CD	PRO	A	626	20.565	88.780	34.858	1.00	13.83	6
4841	N	LEU	A	627	18.725	91.061	31.332	1.00	13.98	7
4842	CA	LEU	A	627	18.935	91.116	29.881	1.00	13.16	6
4843	C	LEU	A	627	20.364	91.609	29.649	1.00	14.14	6
4844	O	LEU	A	627	21.024	92.215	30.527	1.00	14.86	8
4845	CB	LEU	A	627	17.906	92.080	29.214	1.00	13.73	6
4846	CG	LEU	A	627	16.549	91.433	28.919	1.00	13.68	6
4847	CD1	LEU	A	627	15.771	91.075	30.200	1.00	16.08	6
4848	CD2	LEU	A	627	15.630	92.370	28.084	1.00	16.33	7
4849	N	LEU	A	628	20.804	91.374	28.421	1.00	12.55	6
4850	CA	LEU	A	628	22.139	91.805	27.919	1.00	12.76	6
4851	C	LEU	A	628	21.994	92.999	26.973	1.00	15.67	8
4852	O	LEU	A	628	20.910	93.197	26.394	1.00	14.76	6
4853	CB	LEU	A	628	22.746	90.663	27.111	1.00	14.06	6
4854	CG	LEU	A	628	22.848	89.332	27.886	1.00	14.72	6
4855	CD1	LEU	A	628	23.526	89.239	27.056	1.00	12.14	6
4856	CD2	LEU	A	628	23.659	89.508	29.185	1.00	20.12	7
4857	N	ALA	A	629	23.117	93.699	26.702	1.00	16.33	7
4858	CA	ALA	A	629	22.871	94.875	25.787	1.00	16.83	6
4859	C	ALA	A	629	23.983	95.083	24.795	1.00	15.77	6
4860	O	ALA	A	629	24.610	96.177	24.707	1.00	15.92	8
4861	CB	ALA	A	629	22.674	96.098	26.674	1.00	19.56	6
4862	N	PRO	A	630	24.278	94.136	23.970	1.00	19.28	7
4863	CA	PRO	A	630	25.237	94.281	22.873	1.00	19.28	6
4864	C	PRO	A	630	24.773	95.400	21.940	1.00	21.06	6
4865	O	PRO	A	630	25.633	96.015	21.314	1.00	23.58	8
4866	CB	PRO	A	630	25.314	92.944	22.123	1.00	17.59	6
4867	CG	PRO	A	630	23.928	92.367	22.442	1.00	18.42	6
4868	CD	PRO	A	630	23.644	92.807	23.877	1.00	18.05	6
4869	N	ASN	A	631	23.464	95.562	21.768	1.00	17.93	7
4870	CA	ASN	A	631	22.935	96.635	20.926	1.00	17.58	6
4871	C	ASN	A	631	22.382	97.779	21.748	1.00	17.45	6
4872	O	ASN	A	631	21.359	98.370	21.354	1.00	18.41	8
4873	CB	ASN	A	631	21.902	96.082	19.950	1.00	18.61	6
4874	CG	ASN	A	631	22.400	94.876	19.161	1.00	28.21	6
4875	OD1	ASN	A	631	21.838	93.764	19.168	1.00	27.73	8
4876	ND2	ASN	A	631	23.476	95.224	18.455	1.00	19.54	7
4877	N	TYR	A	632	22.980	98.110	22.883	1.00	16.27	7
4878	CA	TYR	A	632	22.570	99.229	23.700	1.00	16.84	6
4879	C	TYR	A	632	22.255	100.496	22.909	1.00	19.89	6
4880	O	TYR	A	632	23.030	100.672	21.982	1.00	19.89	8
4881	CB	TYR	A	632	23.749	99.537	24.643	1.00	18.23	6
4882	CG	TYR	A	632	23.520	100.604	25.654	1.00	20.78	6
4883	CD1	TYR	A	632	23.839	100.369	26.875	1.00	20.07	6
4884	CD2	TYR	A	632	22.919	101.933	25.333	1.00	22.98	6
4885	CE1	TYR	A	632	22.700	101.367	27.797	1.00	22.62	6
4886	CE2	TYR	A	632	23.578	102.953	26.223	1.00	23.64	6
4887	CZ	TYR	A	632	23.051	102.671	27.457	1.00	27.32	6

Serial	Atom	Res	Chain	ResNum	X	Y	Z	Occ	B	N
4888	OH	TYR	A	632	22.814	103.670	28.369	1.00	23.27	8
4889	N	PRO	A	633	21.113	101.067	23.157	1.00	19.06	7
4890	CA	PRO	A	633	20.194	101.103	24.225	1.00	16.58	6
4891	C	PRO	A	633	19.087	100.033	24.123	1.00	15.76	6
4892	O	PRO	A	633	18.188	100.153	24.954	1.00	16.24	8
4893	CB	PRO	A	633	19.474	102.498	24.311	1.00	19.94	6
4894	CG	PRO	A	633	19.449	102.810	22.846	1.00	24.00	6
4895	CD	PRO	A	633	20.772	102.271	22.369	1.00	24.76	6
4896	N	ASP	A	634	19.272	99.094	23.169	1.00	16.18	7
4897	CA	ASP	A	634	18.330	97.978	23.121	1.00	16.48	6
4898	C	ASP	A	634	18.923	96.792	23.937	1.00	15.97	6
4899	O	ASP	A	634	20.153	96.671	24.058	1.00	16.27	8
4900	CB	ASP	A	634	18.133	97.402	21.731	1.00	15.05	6
4901	CG	ASP	A	634	17.626	98.434	20.695	1.00	20.07	6
4902	OD1	ASP	A	634	17.138	99.519	21.100	1.00	17.34	8
4903	OD2	ASP	A	634	17.728	98.097	19.503	1.00	19.32	8
4904	N	TRP	A	635	18.009	96.147	24.656	1.00	14.45	7
4905	CA	TRP	A	635	18.413	95.017	25.510	1.00	16.15	6
4906	C	TRP	A	635	17.708	93.736	25.082	1.00	13.84	6
4907	O	TRP	A	635	16.590	93.820	24.606	1.00	13.90	8
4908	CB	TRP	A	635	18.024	95.299	26.962	1.00	15.52	6
4909	CG	TRP	A	635	18.818	96.421	27.587	1.00	14.07	6
4910	CD1	TRP	A	635	18.737	97.739	27.178	1.00	14.13	6
4911	CD2	TRP	A	635	19.713	96.407	28.694	1.00	16.17	6
4912	NE1	TRP	A	635	19.561	98.533	27.989	1.00	15.61	7
4913	CE2	TRP	A	635	20.179	97.706	28.908	1.00	17.77	6
4914	CE3	TRP	A	635	20.179	95.350	29.511	1.00	18.51	6
4915	CZ2	TRP	A	635	21.071	98.018	29.935	1.00	19.26	6
4916	CZ3	TRP	A	635	21.100	95.671	30.515	1.00	23.17	6
4917	CH2	TRP	A	635	21.514	96.992	30.735	1.00	20.92	7
4918	N	PHE	A	636	18.294	92.561	25.372	1.00	12.41	6
4919	CA	PHE	A	636	17.562	91.335	24.947	1.00	11.57	6
4920	C	PHE	A	636	18.071	90.194	25.855	1.00	12.41	6
4921	O	PHE	A	636	19.204	90.239	26.345	1.00	13.24	8
4922	CB	PHE	A	636	17.761	90.971	23.435	1.00	12.03	6
4923	CG	PHE	A	636	19.030	90.118	23.261	1.00	12.15	6
4924	CD1	PHE	A	636	20.287	90.667	23.292	1.00	17.38	6
4925	CD2	PHE	A	636	18.856	88.751	23.095	1.00	11.02	6
4926	CE1	PHE	A	636	21.417	89.860	23.207	1.00	19.56	6
4927	CE2	PHE	A	636	19.996	87.906	22.997	1.00	12.49	6
4928	CZ	PHE	A	636	21.250	88.466	23.034	1.00	17.21	6
4929	N	TYR	A	637	17.229	89.170	25.966	1.00	12.09	7
4930	CA	TYR	A	637	17.759	87.841	26.425	1.00	12.35	6
4931	C	TYR	A	637	16.714	86.848	25.989	1.00	12.24	6
4932	O	TYR	A	637	15.911	87.135	25.088	1.00	12.49	8
4933	CB	TYR	A	637	18.183	87.770	27.903	1.00	12.77	6
4934	CG	TYR	A	637	19.035	86.513	28.195	1.00	13.33	6
4935	CD1	TYR	A	637	20.253	86.244	27.548	1.00	12.60	6
4936	CD2	TYR	A	637	18.595	85.573	29.127	1.00	10.55	6
4937	CE1	TYR	A	637	21.022	85.134	27.792	1.00	11.85	6
4938	CE2	TYR	A	637	19.321	84.424	29.387	1.00	12.60	6
4939	CZ	TYR	A	637	20.521	84.221	28.728	1.00	11.69	6
4940	OH	TYR	A	637	21.255	83.073	29.016	1.00	11.85	8
4941	N	VAL	A	638	16.857	85.565	26.379	1.00	12.66	7
4942	CA	VAL	A	638	16.033	84.508	25.827	1.00	10.90	6
4943	C	VAL	A	638	15.492	83.693	27.023	1.00	11.04	6
4944	O	VAL	A	638	16.202	83.367	27.980	1.00	13.35	8
4945	CB	VAL	A	638	16.830	83.496	24.987	1.00	12.00	6
4946	CG1	VAL	A	638	15.922	82.650	24.118	1.00	12.74	6
4947	CG2	VAL	A	638	17.799	84.300	24.074	1.00	10.59	6
4948	N	PHE	A	639	14.137	83.517	26.957	1.00	12.77	7
4949	CA	PHE	A	639	13.441	82.960	28.082	1.00	9.34	6
4950	C	PHE	A	639	12.466	81.877	27.665	1.00	11.56	6
4951	O	PHE	A	639	11.814	81.982	26.632	1.00	12.26	8
4952	CB	PHE	A	639	12.612	84.072	28.807	1.00	11.50	6
4953	CG	PHE	A	639	13.498	85.209	29.322	1.00	12.29	6
4954	CD1	PHE	A	639	14.294	85.042	30.446	1.00	11.38	6
4955	CD2	PHE	A	639	13.567	86.380	28.563	1.00	13.67	6
4956	CE1	PHE	A	639	15.176	86.094	30.829	1.00	12.05	6
4957	CE2	PHE	A	639	14.416	87.425	28.965	1.00	13.84	6
4958	CZ	PHE	A	639	15.196	87.301	30.113	1.00	13.28	6
4959	N	SER	A	640	12.323	80.913	28.609	1.00	10.61	7
4960	CA	SER	A	640	11.376	79.833	28.360	1.00	11.72	6
4961	C	SER	A	640	10.005	80.223	28.872	1.00	11.19	6
4962	O	SER	A	640	9.878	80.649	30.026	1.00	12.90	8
4963	CB	SER	A	640	11.915	78.568	29.132	1.00	12.23	6
4964	OG	SER	A	640	11.028	77.448	28.846	1.00	12.04	8
4965	N	VAL	A	641	9.007	80.190	27.975	1.00	8.82	7
4966	CA	VAL	A	641	7.635	80.609	28.374	1.00	11.37	6
4967	C	VAL	A	641	6.697	79.601	27.775	1.00	11.35	6
4968	O	VAL	A	641	7.075	78.840	26.864	1.00	10.31	8
4969	CB	VAL	A	641	7.286	82.019	27.760	1.00	11.50	6
4970	CG1	VAL	A	641	8.061	83.038	28.624	1.00	13.20	6
4971	CG2	VAL	A	641	7.607	82.115	26.241	1.00	10.82	6

No.	Atom	Res	Ch	ResNo	X	Y	Z	Occ	B	El
4972	N	PRO	A	642	5.470	79.384	28.254	1.00	10.92	7
4973	CA	PRO	A	642	4.582	78.438	27.626	1.00	11.56	6
4974	C	PRO	A	642	4.256	78.780	26.166	1.00	10.87	6
4975	O	PRO	A	642	4.048	79.994	25.888	1.00	12.26	8
4976	CB	PRO	A	642	3.295	78.524	28.529	1.00	11.37	6
4977	CG	PRO	A	642	3.352	79.948	29.064	1.00	11.52	6
4978	CD	PRO	A	642	4.866	80.234	29.308	1.00	14.12	6
4979	N	ALA	A	643	4.267	77.800	25.291	1.00	11.34	7
4980	CA	ALA	A	643	4.033	77.992	23.839	1.00	12.12	6
4981	C	ALA	A	643	2.531	78.215	23.567	1.00	13.78	6
4982	O	ALA	A	643	1.681	77.650	24.236	1.00	15.31	8
4983	CB	ALA	A	643	4.330	76.657	23.141	1.00	13.71	6
4984	N	GLY	A	644	2.317	79.106	22.623	1.00	14.27	7
4985	CA	GLY	A	644	0.952	79.377	22.150	1.00	14.42	6
4986	C	GLY	A	644	0.065	80.108	23.133	1.00	15.94	6
4987	O	GLY	A	644	-1.154	79.765	23.178	1.00	19.30	8
4988	N	LYS	A	645	0.597	80.884	24.030	1.00	14.60	7
4989	CA	LYS	A	645	-0.171	81.519	25.092	1.00	14.20	6
4990	C	LYS	A	645	-0.060	83.017	24.926	1.00	17.57	6
4991	O	LYS	A	645	1.002	83.579	24.623	1.00	17.47	8
4992	CB	LYS	A	645	0.376	81.126	26.483	1.00	14.57	6
4993	CG	LYS	A	645	0.313	79.611	26.725	1.00	17.45	6
4994	CD	LYS	A	645	-1.136	79.116	26.699	1.00	17.54	6
4995	CE	LYS	A	645	-1.184	77.572	26.837	1.00	21.69	6
4996	NZ	LYS	A	645	-0.587	77.121	28.112	1.00	31.91	7
4997	N	THR	A	646	-1.157	83.717	25.357	1.00	14.42	7
4998	CA	THR	A	646	-1.090	85.140	25.434	1.00	14.66	6
4999	C	THR	A	646	-0.662	85.594	26.824	1.00	13.55	6
5000	O	THR	A	646	-1.221	85.157	27.850	1.00	16.55	8
5001	CB	THR	A	646	-2.510	85.720	25.098	1.00	17.00	6
5002	OG1	THR	A	646	-2.824	85.361	23.738	1.00	16.76	8
5003	CG2	THR	A	646	-2.463	87.240	25.195	1.00	18.18	6
5004	N	ILE	A	647	0.360	86.434	26.913	1.00	11.81	7
5005	CA	ILE	A	647	1.001	86.816	28.152	1.00	11.47	6
5006	C	ILE	A	647	1.047	88.357	28.250	1.00	13.84	6
5007	O	ILE	A	647	0.991	89.054	27.201	1.00	14.89	8
5008	CB	ILE	A	647	2.476	86.353	28.278	1.00	13.09	6
5009	CG1	ILE	A	647	3.296	86.770	27.045	1.00	12.50	6
5010	CG2	ILE	A	647	2.410	84.817	28.343	1.00	14.24	6
5011	CD1	ILE	A	647	4.813	86.486	27.247	1.00	17.78	6
5012	N	GLN	A	648	1.150	88.776	29.495	1.00	12.61	7
5013	CA	GLN	A	648	1.402	90.212	29.726	1.00	11.91	6
5014	C	GLN	A	648	2.687	90.407	30.490	1.00	15.15	6
5015	O	GLN	A	648	3.041	89.586	31.338	1.00	15.89	8
5016	CB	GLN	A	648	0.298	90.835	30.611	1.00	13.04	6
5017	CG	GLN	A	648	-1.032	90.704	29.820	1.00	20.15	6
5018	CD	GLN	A	648	-2.040	91.727	30.260	1.00	27.13	6
5019	OE1	GLN	A	648	-1.797	92.691	30.982	1.00	23.22	8
5020	NE2	GLN	A	648	-3.309	91.503	29.847	1.00	28.35	7
5021	N	PHE	A	649	3.446	91.461	30.155	1.00	13.26	7
5022	CA	PHE	A	649	4.767	91.622	30.746	1.00	12.72	6
5023	C	PHE	A	649	5.219	93.094	30.663	1.00	14.64	6
5024	O	PHE	A	649	4.697	93.903	29.862	1.00	14.50	8
5025	CB	PHE	A	649	5.812	90.743	30.051	1.00	12.47	6
5026	CG	PHE	A	649	5.875	90.962	28.544	1.00	13.33	6
5027	CD1	PHE	A	649	4.996	90.400	27.682	1.00	14.84	6
5028	CD2	PHE	A	649	6.861	91.790	28.013	1.00	13.41	6
5029	CE1	PHE	A	649	5.031	90.678	26.303	1.00	15.17	6
5030	CE2	PHE	A	649	6.977	92.043	26.666	1.00	16.16	6
5031	CZ	PHE	A	649	6.057	91.493	25.807	1.00	13.88	6
5032	N	LYS	A	650	6.135	93.413	31.538	1.00	13.99	7
5033	CA	LYS	A	650	6.876	94.683	31.405	1.00	13.18	6
5034	C	LYS	A	650	8.349	94.411	31.797	1.00	15.87	6
5035	O	LYS	A	650	8.696	93.364	32.375	1.00	13.96	8
5036	CB	LYS	A	650	6.358	95.735	32.410	1.00	14.62	6
5037	CG	LYS	A	650	5.004	96.385	31.989	1.00	15.36	6
5038	CD	LYS	A	650	4.922	97.604	32.979	1.00	19.44	6
5039	CE	LYS	A	650	3.841	98.565	32.553	1.00	24.45	6
5040	NZ	LYS	A	650	3.898	99.741	33.515	1.00	18.41	7
5041	N	PHE	A	651	9.172	95.398	31.438	1.00	13.40	7
5042	CA	PHE	A	651	10.571	95.420	31.850	1.00	14.23	6
5043	C	PHE	A	651	10.800	96.474	32.926	1.00	16.48	6
5044	O	PHE	A	651	9.990	97.396	33.132	1.00	17.07	8
5045	CB	PHE	A	651	11.465	95.783	30.638	1.00	12.48	6
5046	CG	PHE	A	651	11.280	94.796	29.512	1.00	15.91	6
5047	CD1	PHE	A	651	10.277	94.943	28.564	1.00	13.76	6
5048	CD2	PHE	A	651	12.127	93.681	29.453	1.00	14.89	6
5049	CE1	PHE	A	651	10.114	94.016	27.568	1.00	13.50	6
5050	CE2	PHE	A	651	11.957	92.759	28.426	1.00	12.48	6
5051	CZ	PHE	A	651	10.944	92.884	27.461	1.00	16.56	6
5052	N	PHE	A	652	11.849	96.240	33.723	1.00	15.18	7
5053	CA	PHE	A	652	12.249	97.256	34.704	1.00	14.29	6
5054	C	PHE	A	652	13.756	97.339	34.785	1.00	17.92	6
5055	O	PHE	A	652	14.452	96.365	34.487	1.00	15.37	8

5056	CB	PHE	A	652	11.634	96.950	36.084	1.00	15.14	6
5057	CG	PHE	A	652	12.097	95.688	36.789	1.00	16.77	6
5058	CD1	PHE	A	652	13.251	95.685	37.565	1.00	15.74	6
5059	CD2	PHE	A	652	11.326	94.553	36.695	1.00	15.90	6
5060	CE1	PHE	A	652	13.633	94.518	38.231	1.00	14.76	6
5061	CE2	PHE	A	652	11.695	93.407	37.380	1.00	15.19	6
5062	CZ	PHE	A	652	12.866	93.343	38.139	1.00	14.25	6
5063	N	ILE	A	653	14.244	98.512	35.189	1.00	14.94	7
5064	CA	ILE	A	653	15.705	98.598	35.463	1.00	14.72	6
5065	C	ILE	A	653	15.906	98.394	36.962	1.00	18.67	6
5066	O	ILE	A	653	15.175	98.971	37.804	1.00	19.59	8
5067	CB	ILE	A	653	16.168	100.071	35.163	1.00	14.73	6
5068	CG1	ILE	A	653	16.000	100.337	33.693	1.00	18.74	6
5069	CG2	ILE	A	653	17.628	100.201	35.616	1.00	18.25	6
5070	CD1	ILE	A	653	16.191	101.838	33.353	1.00	20.43	6
5071	N	LYS	A	654	16.847	97.524	37.282	1.00	17.58	7
5072	CA	LYS	A	654	17.273	97.382	38.690	1.00	18.28	6
5073	C	LYS	A	654	18.623	98.139	38.780	1.00	15.87	6
5074	O	LYS	A	654	19.571	97.760	38.135	1.00	18.77	8
5075	CB	LYS	A	654	17.469	95.933	39.131	1.00	16.77	6
5076	CG	LYS	A	654	17.774	95.976	40.661	1.00	20.67	6
5077	CE	LYS	A	654	17.709	94.558	41.225	1.00	24.12	6
5078	CE	LYS	A	654	18.152	94.544	42.680	1.00	30.47	6
5079	NZ	LYS	A	654	17.967	93.152	43.221	1.00	39.55	7
5080	N	ARG	A	655	18.538	99.361	39.337	1.00	19.37	7
5081	CA	ARG	A	655	19.776	100.164	39.391	1.00	21.00	6
5082	C	ARG	A	655	20.820	99.478	40.256	1.00	24.58	6
5083	O	ARG	A	655	20.486	98.604	41.057	1.00	22.97	8
5084	CB	ARG	A	655	19.445	101.531	40.003	1.00	18.71	6
5085	CG	ARG	A	655	18.411	102.323	39.192	1.00	19.26	6
5086	CD	ARG	A	655	19.036	102.699	37.864	1.00	24.72	6
5087	NE	ARG	A	655	18.169	103.597	37.102	1.00	23.26	7
5088	CZ	ARG	A	655	18.428	104.073	35.883	1.00	23.74	6
5089	NH1	ARG	A	655	19.508	103.812	35.193	1.00	20.63	7
5090	NH2	ARG	A	655	17.509	104.882	35.343	1.00	24.30	7
5091	N	ALA	A	656	22.090	99.900	40.135	1.00	24.51	7
5092	CA	ALA	A	656	23.170	99.305	40.898	1.00	28.24	6
5093	C	ALA	A	656	22.865	99.430	42.401	1.00	27.25	6
5094	O	ALA	A	656	23.305	98.533	43.117	1.00	32.90	8
5095	CB	ALA	A	656	24.518	100.009	40.656	1.00	29.09	6
5096	N	ASP	A	657	22.145	100.427	42.856	1.00	31.34	7
5097	CA	ASP	A	657	21.850	100.561	44.282	1.00	32.95	6
5098	C	ASP	A	657	20.617	99.783	44.694	1.00	33.22	6
5099	O	ASP	A	657	20.202	99.906	45.852	1.00	30.17	8
5100	CB	ASP	A	657	21.725	102.050	44.643	1.00	27.78	6
5101	CG	ASP	A	657	20.499	102.712	44.087	1.00	34.66	6
5102	OD1	ASP	A	657	19.665	102.035	43.431	1.00	28.22	8
5103	OD2	ASP	A	657	20.239	103.920	44.267	1.00	33.52	8
5104	N	GLY	A	658	20.019	98.958	43.803	1.00	29.54	7
5105	CA	GLY	A	658	18.874	98.181	44.227	1.00	29.70	6
5106	C	GLY	A	658	17.543	98.837	43.931	1.00	26.81	6
5107	O	GLY	A	658	16.531	98.126	43.935	1.00	30.39	8
5108	N	THR	A	659	17.502	100.106	43.566	1.00	23.52	7
5109	CA	THR	A	659	16.313	100.826	43.215	1.00	25.61	6
5110	C	THR	A	659	15.641	100.184	41.985	1.00	23.40	6
5111	O	THR	A	659	16.419	99.778	41.118	1.00	27.85	8
5112	CB	THR	A	659	16.614	102.279	42.717	1.00	19.42	6
5113	OG1	THR	A	659	17.524	102.993	43.510	1.00	46.62	8
5114	CG2	THR	A	659	15.276	103.016	42.706	1.00	37.02	6
5115	N	ILE	A	660	14.319	100.215	41.909	1.00	20.34	7
5116	CA	ILE	A	660	13.596	99.691	40.758	1.00	25.29	6
5117	C	ILE	A	660	12.933	100.761	39.919	1.00	25.57	6
5118	O	ILE	A	660	12.186	101.612	40.465	1.00	27.02	8
5119	CB	ILE	A	660	12.458	98.722	41.207	1.00	26.69	6
5120	CG1	ILE	A	660	13.012	97.534	41.991	1.00	27.58	6
5121	CG2	ILE	A	660	11.704	98.273	39.972	1.00	21.18	6
5122	CD1	ILE	A	660	14.022	96.722	41.221	1.00	24.46	6
5123	N	GLN	A	661	13.169	100.849	38.624	1.00	20.86	7
5124	CA	GLN	A	661	12.479	101.793	37.743	1.00	19.49	6
5125	C	GLN	A	661	11.720	101.014	36.669	1.00	20.02	6
5126	O	GLN	A	661	12.266	100.405	35.730	1.00	19.24	8
5127	CB	GLN	A	661	13.410	102.775	37.023	1.00	20.06	6
5128	CG	GLN	A	661	12.611	103.688	36.077	1.00	21.00	6
5129	CD	GLN	A	661	13.541	104.665	35.350	1.00	21.47	6
5130	OE1	GLN	A	661	13.279	104.997	34.195	1.00	28.51	8
5131	NE2	GLN	A	661	14.578	105.095	36.040	1.00	22.67	7
5132	N	TRP	A	662	10.389	100.926	36.834	1.00	20.75	7
5133	CA	TRP	A	662	9.556	100.227	35.857	1.00	18.49	6
5134	C	TRP	A	662	9.441	100.967	34.536	1.00	21.67	6
5135	O	TRP	A	662	9.412	102.230	34.565	1.00	19.29	8
5136	CB	TRP	A	662	8.152	100.082	36.437	1.00	18.81	6
5137	CG	TRP	A	662	7.960	99.114	37.551	1.00	18.39	6
5138	CD1	TRP	A	662	8.083	99.359	38.889	1.00	19.35	6
5139	CD2	TRP	A	662	7.543	97.752	37.419	1.00	18.13	6

5140	NE1	TRP	A	662	7.781	98.218	39.608	1.00	21.46	7
5141	CE2	TRP	A	662	7.444	97.223	38.713	1.00	18.92	6
5142	CE3	TRP	A	662	7.213	96.960	36.295	1.00	18.78	6
5143	CZ2	TRP	A	662	6.994	95.923	38.966	1.00	16.39	6
5144	CZ3	TRP	A	662	6.798	95.661	36.537	1.00	18.67	6
5145	CH2	TRP	A	662	6.725	95.156	37.854	1.00	15.44	6
5146	N	GLU	A	663	9.170	100.250	33.453	1.00	18.24	7
5147	CA	GLU	A	663	8.551	100.857	32.292	1.00	17.79	6
5148	C	GLU	A	663	7.240	101.517	32.740	1.00	17.95	6
5149	O	GLU	A	663	6.519	100.973	33.574	1.00	18.49	8
5150	CB	GLU	A	663	8.001	99.880	31.232	1.00	22.54	6
5151	CG	GLU	A	663	9.065	99.526	30.195	1.00	23.03	6
5152	CD	GLU	A	663	8.380	98.560	29.223	1.00	19.52	6
5153	OE1	GLU	A	663	8.159	97.397	29.619	1.00	17.12	8
5154	OE2	GLU	A	663	8.063	98.990	28.102	1.00	18.01	8
5155	N	ASN	A	664	6.892	102.576	32.027	1.00	18.62	7
5156	CA	ASN	A	664	5.604	103.189	32.365	1.00	18.27	6
5157	C	ASN	A	664	4.522	102.581	31.484	1.00	20.85	6
5158	O	ASN	A	664	4.782	101.671	30.702	1.00	21.34	8
5159	CB	ASN	A	664	5.790	104.696	32.115	1.00	22.20	6
5160	CG	ASN	A	664	6.447	105.326	33.350	1.00	30.00	6
5161	OD1	ASN	A	664	6.281	104.932	34.519	1.00	35.11	8
5162	ND2	ASN	A	664	7.238	106.366	33.127	1.00	33.46	7
5163	N	GLY	A	665	3.295	103.085	31.667	1.00	22.07	6
5164	CA	GLY	A	665	2.187	102.708	30.786	1.00	23.10	6
5165	C	GLY	A	665	1.557	101.363	31.102	1.00	21.34	6
5166	O	GLY	A	665	1.827	100.710	32.119	1.00	22.70	8
5167	N	SER	A	666	0.832	100.839	30.093	1.00	18.43	7
5168	CA	SER	A	666	0.135	99.581	30.305	1.00	17.79	6
5169	C	SER	A	666	1.084	98.403	30.046	1.00	16.94	6
5170	O	SER	A	666	2.005	98.503	29.261	1.00	16.68	8
5171	CB	SER	A	666	-1.037	99.440	29.297	1.00	24.19	6
5172	OG	SER	A	666	-1.959	100.498	29.645	1.00	24.86	8
5173	N	ASN	A	667	0.642	97.227	30.470	1.00	14.27	7
5174	CA	ASN	A	667	1.491	96.058	30.168	1.00	16.75	6
5175	C	ASN	A	667	1.575	95.824	28.678	1.00	17.00	6
5176	O	ASN	A	667	0.616	96.050	27.899	1.00	15.91	8
5177	CB	ASN	A	667	0.795	94.804	30.736	1.00	16.66	6
5178	CG	ASN	A	667	0.703	94.817	32.248	1.00	19.33	6
5179	OD1	ASN	A	667	-0.178	94.105	32.855	1.00	20.96	8
5180	ND2	ASN	A	667	1.594	95.513	32.887	1.00	16.21	7
5181	N	HIS	A	668	2.720	95.243	28.272	1.00	13.63	7
5182	CA	HIS	A	668	2.882	94.747	26.930	1.00	15.66	6
5183	C	HIS	A	668	2.137	93.408	26.856	1.00	16.10	6
5184	O	HIS	A	668	2.048	92.666	27.847	1.00	16.87	8
5185	CB	HIS	A	668	4.327	94.431	26.491	1.00	15.51	6
5186	CG	HIS	A	668	5.208	95.653	26.519	1.00	14.13	6
5187	ND1	HIS	A	668	5.229	96.451	25.375	1.00	15.75	6
5188	CD2	HIS	A	668	6.066	96.158	27.410	1.00	17.00	7
5189	CE1	HIS	A	668	6.108	97.464	25.592	1.00	14.21	6
5190	NE2	HIS	A	668	6.581	97.305	26.830	1.00	15.51	6
5191	N	VAL	A	669	1.434	93.208	25.734	1.00	15.02	7
5192	CA	VAL	A	669	0.684	91.987	25.490	1.00	16.94	7
5193	C	VAL	A	669	1.205	91.294	24.260	1.00	14.25	6
5194	O	VAL	A	669	1.414	91.936	23.206	1.00	17.08	8
5195	CB	VAL	A	669	-0.852	92.295	25.322	1.00	17.60	6
5196	CG1	VAL	A	669	-1.624	90.970	25.253	1.00	20.63	6
5197	CG2	VAL	A	669	-1.341	93.078	26.547	1.00	17.53	6
5198	N	ALA	A	670	1.450	89.971	24.287	1.00	14.50	7
5199	CA	ALA	A	670	1.945	89.249	23.113	1.00	14.21	6
5200	C	ALA	A	670	1.437	87.798	23.234	1.00	15.36	8
5201	O	ALA	A	670	1.216	87.262	24.323	1.00	15.04	6
5202	CB	ALA	A	670	3.481	89.251	23.005	1.00	14.71	6
5203	N	THR	A	671	3.412	87.143	22.093	1.00	15.66	8
5204	CA	THR	A	671	1.145	85.704	22.056	1.00	15.06	6
5205	C	THR	A	671	2.463	85.017	21.737	1.00	15.13	7
5206	O	THR	A	671	3.141	85.293	20.734	1.00	16.78	6
5207	CB	THR	A	671	0.004	85.374	21.082	1.00	24.32	6
5208	OG1	THR	A	671	-1.181	86.041	21.584	1.00	19.33	8
5209	CG2	THR	A	671	-0.323	83.900	21.063	1.00	22.47	7
5210	N	THR	A	672	2.794	83.967	22.533	1.00	13.36	6
5211	CA	THR	A	672	4.035	83.264	22.249	1.00	14.54	6
5212	C	THR	A	672	3.911	82.312	21.096	1.00	14.80	8
5213	O	THR	A	672	2.831	81.798	20.772	1.00	13.92	6
5214	CB	THR	A	672	4.475	82.464	23.491	1.00	15.06	8
5215	OG1	THR	A	672	3.485	81.497	23.805	1.00	13.92	6
5216	CG2	THR	A	672	4.612	83.376	24.726	1.00	15.62	6
5217	N	PRO	A	673	5.068	81.983	20.499	1.00	16.20	6
5218	CA	PRO	A	673	5.116	81.063	19.381	1.00	21.02	8
5219	C	PRO	A	673	4.615	79.690	19.691	1.00	17.09	6
5220	O	PRO	A	673	4.691	79.290	20.893	1.00	16.99	8
5221	CB	PRO	A	673	6.606	80.975	18.937	1.00	22.47	6
5222	CG	PRO	A	673	7.309	81.859	19.872	1.00	22.50	6
5223	CD	PRO	A	673	6.355	82.558	20.857	1.00	20.76	6

5224	N	THR A 674	4.164	78.885	18.724	1.00	16.36	7
5225	CA	THR A 674	3.803	77.518	19.033	1.00	16.04	6
5226	C	THR A 674	4.915	76.516	18.803	1.00	19.50	6
5227	O	THR A 674	4.834	75.399	19.293	1.00	25.45	8
5228	CB	THR A 674	2.613	77.053	18.117	1.00	26.38	6
5229	OG1	THR A 674	2.997	77.353	16.786	1.00	31.27	8
5230	CG2	THR A 674	1.409	77.936	18.521	1.00	27.63	6
5231	N	GLY A 675	5.953	76.981	18.114	1.00	19.68	7
5232	CA	GLY A 675	7.035	76.041	17.829	1.00	18.98	6
5233	C	GLY A 675	8.164	76.145	18.885	1.00	21.20	6
5234	O	GLY A 675	7.915	76.654	19.953	1.00	18.10	8
5235	N	ALA A 676	9.349	75.612	18.560	1.00	14.92	7
5236	CA	ALA A 676	10.419	75.562	19.578	1.00	12.67	6
5237	C	ALA A 676	10.845	76.968	19.974	1.00	13.87	6
5238	O	ALA A 676	11.505	77.096	21.031	1.00	12.93	8
5239	CB	ALA A 676	11.611	74.771	19.006	1.00	12.85	6
5240	N	THR A 677	10.860	77.884	19.014	1.00	13.69	7
5241	CA	THR A 677	11.444	79.189	19.324	1.00	11.81	6
5242	C	THR A 677	10.615	80.296	18.686	1.00	13.20	6
5243	O	THR A 677	9.743	80.066	17.834	1.00	14.64	8
5244	CB	THR A 677	12.847	79.366	18.681	1.00	14.32	6
5245	OG1	THR A 677	12.712	79.294	17.249	1.00	16.35	8
5246	CG2	THR A 677	13.851	78.287	19.079	1.00	14.03	6
5247	N	GLY A 678	10.949	81.490	19.141	1.00	14.25	7
5248	CA	GLY A 678	10.413	82.685	18.437	1.00	13.38	6
5249	C	GLY A 678	10.969	83.958	19.066	1.00	13.68	6
5250	O	GLY A 678	11.857	83.956	19.907	1.00	13.62	8
5251	N	ASN A 679	10.374	85.094	18.654	1.00	13.18	7
5252	CA	ASN A 679	10.858	86.398	19.082	1.00	13.18	6
5253	C	ASN A 679	9.685	87.318	19.466	1.00	11.86	6
5254	O	ASN A 679	8.718	87.315	18.699	1.00	14.28	8
5255	CB	ASN A 679	11.485	87.180	17.892	1.00	15.54	6
5256	CG	ASN A 679	12.952	86.818	17.758	1.00	16.45	6
5257	OD1	ASN A 679	13.201	85.708	17.294	1.00	25.76	8
5258	ND2	ASN A 679	13.881	87.684	18.146	1.00	24.19	7
5259	N	ILE A 680	9.882	88.063	20.521	1.00	26.44	7
5260	CA	ILE A 680	8.909	89.088	20.948	1.00	13.28	6
5261	C	ILE A 680	9.758	90.331	21.083	1.00	14.10	6
5262	O	ILE A 680	10.673	90.366	21.944	1.00	16.54	8
5263	CB	ILE A 680	8.224	88.676	22.277	1.00	14.10	6
5264	CG1	ILE A 680	7.245	87.519	21.978	1.00	16.92	6
5265	CG2	ILE A 680	7.442	89.919	22.790	1.00	17.42	6
5266	CD1	ILE A 680	6.611	86.867	23.218	1.00	15.15	6
5267	N	THR A 681	9.363	91.378	20.334	1.00	13.57	7
5268	CA	THR A 681	10.236	92.583	20.392	1.00	12.91	6
5269	C	THR A 681	9.280	93.751	20.616	1.00	13.01	6
5270	O	THR A 681	8.253	93.845	19.929	1.00	18.59	8
5271	CB	THR A 681	10.904	92.718	18.990	1.00	16.97	6
5272	OG1	THR A 681	11.807	91.633	18.765	1.00	17.30	8
5273	CG2	THR A 681	11.642	94.052	18.964	1.00	19.53	6
5274	N	VAL A 682	9.639	94.595	21.608	1.00	15.51	7
5275	CA	VAL A 682	8.758	95.732	21.946	1.00	15.78	6
5276	C	VAL A 682	9.635	96.952	22.186	1.00	15.87	6
5277	O	VAL A 682	10.838	96.797	22.256	1.00	15.49	8
5278	CB	VAL A 682	7.874	95.472	23.195	1.00	16.15	6
5279	CG1	VAL A 682	6.968	94.243	23.002	1.00	16.76	6
5280	CG2	VAL A 682	8.730	95.365	24.469	1.00	16.17	8
5281	N	THR A 683	9.082	98.162	22.306	1.00	15.89	6
5282	CA	THR A 683	9.863	99.346	22.623	1.00	15.51	6
5283	C	THR A 683	9.529	99.798	24.046	1.00	16.36	6
5284	O	THR A 683	8.371	99.703	24.462	1.00	17.59	8
5285	CB	THR A 683	9.481	100.473	21.614	1.00	22.31	8
5286	OG1	THR A 683	9.916	100.019	20.328	1.00	21.22	8
5287	CG2	THR A 683	10.245	101.759	21.921	1.00	19.31	6
5288	N	TRP A 684	10.586	100.168	24.759	1.00	14.24	7
5289	CA	TRP A 684	10.411	100.587	26.154	1.00	14.76	6
5290	C	TRP A 684	9.369	101.706	26.222	1.00	17.75	6
5291	O	TRP A 684	9.556	102.695	25.469	1.00	18.74	8
5292	CB	TRP A 684	11.745	101.086	26.705	1.00	14.95	6
5293	CG	TRP A 684	11.696	101.465	28.146	1.00	14.49	6
5294	CD1	TRP A 684	11.284	102.666	28.672	1.00	15.39	6
5295	CD2	TRP A 684	12.163	100.683	29.271	1.00	17.84	6
5296	NE1	TRP A 684	11.376	102.649	30.052	1.00	16.11	7
5297	CE2	TRP A 684	11.921	101.440	30.423	1.00	20.66	6
5298	CE3	TRP A 684	12.764	99.422	29.399	1.00	18.28	6
5299	CZ2	TRP A 684	12.245	101.024	31.726	1.00	22.23	8
5300	CZ3	TRP A 684	13.074	98.959	30.709	1.00	17.05	6
5301	CH2	TRP A 684	12.791	99.766	31.796	1.00	17.97	6
5302	N	GLN A 685	8.396	101.650	27.102	1.00	20.66	7
5303	CA	GLN A 685	7.399	102.716	27.315	1.00	17.70	6
5304	C	GLN A 685	7.850	103.719	28.333	1.00	18.97	6
5305	O	GLN A 685	8.067	103.430	29.512	1.00	19.08	8
5306	CB	GLN A 685	6.071	102.025	27.745	1.00	15.69	6
5307	CG	GLN A 685	5.536	101.157	26.596	1.00	17.89	6

Serial	Atom	Res	Chain	ResNum	X	Y	Z	Occ	B	Elem
5308	CD	GLN	A	685	4.352	100.283	26.995	1.00	18.78	6
5309	OE1	GLN	A	685	3.797	99.604	26.110	1.00	20.58	8
5310	NE2	GLN	A	685	3.960	100.246	28.267	1.00	17.75	7
5311	N	ASN	A	686	8.037	105.011	27.906	1.00	21.27	7
5312	CA	ASN	A	686	8.547	106.008	28.836	1.00	20.38	6
5313	C	ASN	A	686	7.489	106.854	29.654	1.00	22.87	6
5314	O	ASN	A	686	6.387	106.692	29.016	1.00	22.87	8
5315	CB	ASN	A	686	9.300	107.060	27.975	1.00	24.12	6
5316	CG	ASN	A	686	10.434	106.404	27.176	1.00	23.38	6
5317	OD1	ASN	A	686	11.360	105.896	27.821	1.00	22.76	8
5318	ND2	ASN	A	686	10.376	106.412	25.832	1.00	25.47	7
5319	C11	HEX	A	690	38.644	78.012	38.228	1.00	27.46	6
5320	O11	HEX	A	690	39.147	78.243	39.503	1.00	37.24	8
5321	C12	HEX	A	690	37.596	79.172	38.058	1.00	23.95	6
5322	O12	HEX	A	690	36.681	78.887	39.109	1.00	20.96	8
5323	C13	HEX	A	690	36.915	78.887	36.685	1.00	17.64	6
5324	O13	HEX	A	690	35.915	79.908	36.442	1.00	18.04	8
5325	C14	HEX	A	690	38.048	79.023	35.670	1.00	17.74	6
5326	C15	HEX	A	690	39.141	77.967	35.956	1.00	19.52	6
5327	O15	HEX	A	690	39.679	78.312	37.276	1.00	26.20	8
5328	C16	HEX	A	690	40.338	77.919	35.016	1.00	23.71	6
5329	O16	HEX	A	690	40.867	79.240	34.817	1.00	28.38	8
5330	C21	HEX	A	690	37.609	79.324	33.250	1.00	19.03	6
5331	O21	HEX	A	690	37.414	78.562	34.414	1.00	18.39	8
5332	C22	HEX	A	690	36.237	79.756	32.692	1.00	18.97	6
5333	O22	HEX	A	690	35.419	80.404	33.669	1.00	17.96	8
5334	C23	HEX	A	690	35.514	78.516	32.153	1.00	16.72	6
5335	O23	HEX	A	690	34.355	79.014	31.406	1.00	16.31	8
5336	C24	HEX	A	690	36.415	77.760	31.174	1.00	12.77	6
5337	C25	HEX	A	690	37.690	77.301	31.955	1.00	17.32	6
5338	O25	HEX	A	690	38.321	78.596	32.286	1.00	20.72	8
5339	C26	HEX	A	690	38.704	76.681	30.979	1.00	18.94	6
5340	O26	HEX	A	690	39.852	76.182	31.780	1.00	21.07	8
5341	C31	HEX	A	690	35.507	76.403	29.409	1.00	11.28	6
5342	O31	HEX	A	690	35.723	76.501	30.811	1.00	12.85	8
5343	C32	HEX	A	690	34.026	76.118	29.118	1.00	13.04	6
5344	O32	HEX	A	690	33.259	77.030	29.933	1.00	14.06	8
5345	C33	HEX	A	690	33.654	74.698	29.626	1.00	13.59	6
5346	O33	HEX	A	690	32.285	74.486	29.158	1.00	12.48	8
5347	C34	HEX	A	690	34.578	73.670	28.936	1.00	11.83	6
5348	C35	HEX	A	690	36.004	74.036	29.487	1.00	14.01	6
5349	O35	HEX	A	690	36.301	75.341	28.930	1.00	12.70	8
5350	C36	HEX	A	690	37.084	73.167	28.829	1.00	13.71	6
5351	C41	HEX	A	690	33.322	71.601	28.292	1.00	9.18	6
5352	N41	HEX	A	690	34.214	72.300	29.267	1.00	10.73	7
5353	C42	HEX	A	690	33.682	70.074	28.156	1.00	11.48	6
5354	O42	HEX	A	690	33.732	69.433	29.448	1.00	10.30	8
5355	C43	HEX	A	690	35.023	69.963	27.455	1.00	10.39	6
5356	O43	HEX	A	690	35.686	68.699	27.591	1.00	11.44	8
5357	C44	HEX	A	690	34.791	70.257	25.986	1.00	11.02	6
5358	C45	HEX	A	690	33.925	71.520	25.899	1.00	10.58	6
5359	C40	HEX	A	690	33.262	72.097	26.944	1.00	9.08	6
5360	C46	HEX	A	690	33.519	71.887	24.466	1.00	12.28	6
5361	O46	HEX	A	690	32.492	71.022	24.020	1.00	12.58	8
5362	C51	HEX	A	690	36.288	69.864	24.116	1.00	10.88	6
5363	O51	HEX	A	690	36.061	70.541	25.327	1.00	12.10	8
5364	C52	HEX	A	690	37.495	68.906	24.274	1.00	11.69	6
5365	O52	HEX	A	690	37.227	67.917	25.285	1.00	12.57	8
5366	C53	HEX	A	690	38.717	69.698	24.774	1.00	10.51	6
5367	O53	HEX	A	690	39.832	68.756	24.609	1.00	11.97	8
5368	C54	HEX	A	690	39.025	70.807	23.714	1.00	12.82	6
5369	C55	HEX	A	690	37.748	71.660	23.566	1.00	12.56	6
5370	O55	HEX	A	690	36.680	70.820	23.133	1.00	11.57	8
5371	C56	HEX	A	690	37.890	72.687	22.436	1.00	12.61	6
5372	O56	HEX	A	690	38.082	72.074	21.134	1.00	19.74	8
5373	C61	HEX	A	690	40.720	72.590	24.188	1.00	19.96	6
5374	O61	HEX	A	690	39.890	71.572	24.637	1.00	20.26	8
5375	C62	HEX	A	690	42.050	72.500	24.991	1.00	23.13	6
5376	O62	HEX	A	690	42.582	71.189	24.918	1.00	23.06	8
5377	C63	HEX	A	690	41.937	73.006	26.411	1.00	24.03	6
5378	O63	HEX	A	690	43.280	72.994	26.932	1.00	28.78	8
5379	C64	HEX	A	690	41.310	74.405	26.428	1.00	25.79	6
5380	O64	HEX	A	690	41.035	74.789	27.809	1.00	29.96	8
5381	C65	HEX	A	690	39.918	74.255	25.759	1.00	22.19	6
5382	O65	HEX	A	690	40.146	73.866	24.404	1.00	22.47	8
5383	C66	HEX	A	690	39.177	75.555	25.611	1.00	26.12	6
5384	O66	HEX	A	690	39.936	76.644	25.149	1.00	24.41	8
5385	C11	MAL	A	691	38.534	71.299	69.464	1.00	19.75	6
5386	O11	MAL	A	691	38.776	72.581	69.883	1.00	20.21	8
5387	C12	MAL	A	691	37.973	71.384	68.024	1.00	17.66	6
5388	O12	MAL	A	691	38.798	72.262	67.225	1.00	18.71	8
5389	C13	MAL	A	691	36.555	71.949	68.006	1.00	16.03	6
5390	O13	MAL	A	691	35.969	71.754	66.689	1.00	17.47	8
5391	C14	MAL	A	691	35.642	71.104	68.925	1.00	15.64	6

5392	C15	MAL A 691	36.269	71.291	70.364	1.00	14.01	6
5393	O15	MAL A 691	37.561	70.667	70.312	1.00	16.48	8
5394	C16	MAL A 691	35.519	70.221	71.238	1.00	19.77	6
5395	O16	MAL A 691	36.004	70.431	72.581	1.00	18.54	8
5396	C21	MAL A 691	33.285	70.813	68.523	1.00	18.83	6
5397	O21	MAL A 691	34.336	71.683	68.968	1.00	17.92	8
5398	C22	MAL A 691	32.403	71.638	67.561	1.00	17.24	6
5399	O22	MAL A 691	33.177	72.083	66.433	1.00	17.89	8
5400	C23	MAL A 691	31.765	72.820	68.304	1.00	18.41	6
5401	O23	MAL A 691	30.812	73.419	67.431	1.00	18.30	8
5402	C24	MAL A 691	30.951	72.194	69.478	1.00	18.57	6
5403	O24	MAL A 691	30.444	73.291	70.263	1.00	19.15	8
5404	C25	MAL A 691	31.923	71.424	70.383	1.00	19.32	6
5405	O25	MAL A 691	32.521	70.374	69.608	1.00	18.98	8
5406	C26	MAL A 691	31.067	70.708	71.468	1.00	15.38	6
5407	O26	MAL A 691	31.944	70.075	72.412	1.00	18.74	8
5411	S	SUL A 695	11.120	50.018	55.465	1.00	30.54	16
5412	O1	SUL A 695	11.470	52.936	56.533	1.00	30.07	8
5413	O2	SUL A 695	10.034	52.528	54.544	1.00	27.19	8
5414	O3	SUL A 695	12.310	51.631	54.662	1.00	34.25	8
5415	O4	SUL A 695	10.566	50.749	56.089	1.00	33.96	8
5451	CA	WAT A 692	32.693	60.307	13.017	1.00	11.99	20
5452	CA	WAT A 693	26.975	79.502	21.970	1.00	10.73	20
5453	CA	WAT A 694	37.244	49.841	19.039	1.00	13.50	20
5454	OWO	WAT V 1	24.447	79.971	21.858	1.00	9.43	8
5455	OWO	WAT V 2	35.686	59.385	24.028	1.00	10.45	8
5456	OWO	WAT V 3	33.934	60.773	18.648	1.00	10.63	8
5457	OWO	WAT V 4	35.622	62.751	41.495	1.00	10.88	8
5458	OWO	WAT V 5	25.780	77.914	20.486	1.00	10.75	8
5459	OWO	WAT V 6	21.776	77.285	28.879	1.00	10.81	8
5460	OWO	WAT V 7	29.415	69.145	19.400	1.00	10.86	8
5461	OWO	WAT V 8	29.138	80.312	22.631	1.00	10.86	8
5462	OWO	WAT V 9	27.613	72.037	24.448	1.00	10.77	8
5463	OWO	WAT V 10	31.164	77.784	19.615	1.00	10.85	8
5464	OWO	WAT V 11	32.790	66.917	36.378	1.00	10.94	8
5465	OWO	WAT V 12	34.127	66.240	32.929	1.00	11.18	8
5466	OWO	WAT V 13	33.080	60.581	23.767	1.00	11.11	8
5467	OWO	WAT V 14	37.235	54.601	19.600	1.00	11.38	8
5468	OWO	WAT V 15	26.119	65.542	21.574	1.00	11.07	8
5469	OWO	WAT V 16	28.484	64.486	20.522	1.00	11.49	8
5470	OWO	WAT V 17	28.194	73.536	36.853	1.00	11.53	8
5471	OWO	WAT V 18	16.618	66.275	36.806	1.00	11.68	8

5472	OWO	WAT V 19	26.088	62.658	22.854	1.00	11.54	8
5473	OWO	WAT V 20	37.981	63.919	14.127	1.00	11.77	8
5474	OWO	WAT V 21	34.932	60.656	13.769	1.00	11.70	8
5475	OWO	WAT V 22	41.499	60.656	38.722	1.00	11.54	8
5476	OWO	WAT V 23	40.945	66.711	20.205	1.00	11.65	8
5477	OWO	WAT V 24	8.905	64.107	34.370	1.00	11.68	8
5478	OWO	WAT V 25	19.426	72.356	40.893	1.00	11.74	8
5479	OWO	WAT V 26	20.321	82.331	35.376	1.00	11.77	8
5480	OWO	WAT V 27	14.993	64.250	37.502	1.00	11.91	8
5481	OWO	WAT V 28	31.504	68.673	10.842	1.00	12.08	8
5482	OWO	WAT V 29	37.606	61.402	40.167	1.00	11.91	8
5483	OWO	WAT V 30	16.372	70.863	38.933	1.00	11.99	8
5484	OWO	WAT V 31	7.950	69.079	31.258	1.00	12.08	8
5485	OWO	WAT V 32	19.528	73.999	43.164	1.00	12.10	8
5486	OWO	WAT V 33	16.210	66.954	39.606	1.00	12.02	8
5487	OWO	WAT V 34	32.679	63.267	29.330	1.00	12.26	8
5488	OWO	WAT V 35	13.649	74.479	39.683	1.00	12.21	8
5489	OWO	WAT V 36	16.357	79.631	22.013	1.00	12.26	8
5490	OWO	WAT V 37	21.471	63.888	43.225	1.00	12.25	8
5491	OWO	WAT V 38	42.464	66.587	23.881	1.00	12.49	8
5492	OWO	WAT V 39	31.355	60.893	19.838	1.00	12.31	8
5493	OWO	WAT V 40	16.930	64.546	40.981	1.00	12.36	8
5494	OWO	WAT V 41	10.918	81.011	32.425	1.00	12.42	8
5495	OWO	WAT V 42	8.358	78.559	35.685	1.00	12.49	8
5496	OWO	WAT V 43	22.052	71.621	41.190	1.00	12.24	8
5497	OWO	WAT V 44	8.226	66.640	35.276	1.00	12.62	8
5498	OWO	WAT V 45	6.031	77.562	36.868	1.00	12.35	8
5499	OWO	WAT V 46	43.919	60.734	40.175	1.00	12.51	8
5500	OWO	WAT V 48	11.578	73.478	41.191	1.00	12.62	8
5501	OWO	WAT V 49	35.256	52.308	26.203	1.00	12.50	8
5502	OWO	WAT V 50	22.628	81.739	22.782	1.00	12.52	8
5503	OWO	WAT V 51	41.171	68.357	22.292	1.00	12.78	8
5504	OWO	WAT V 52	34.554	71.110	10.783	1.00	12.58	8
5505	OWO	WAT V 53	39.554	70.543	28.407	1.00	12.73	8
5506	OWO	WAT V 54	14.970	66.671	43.779	1.00	12.94	8
5507	OWO	WAT V 55	14.792	81.581	20.763	1.00	12.92	8
5508	OWO	WAT V 56	30.205	75.643	31.109	1.00	12.90	8
5509	OWO	WAT V 57	16.697	82.536	30.534	1.00	13.09	8
5510	OWO	WAT V 58	38.776	53.289	57.732	1.00	13.10	8
5511	OWO	WAT V 59	31.565	49.273	20.064	1.00	13.22	8
5512	OWO	WAT V 60	20.200	85.147	36.037	1.00	13.29	8
5513	OWO	WAT V 61	25.657	54.513	45.949	1.00	13.32	8

ID								
5514	OWO	WAT	V 62	37.048	52.273	18.284	1.00	13.46	8
5515	OWO	WAT	V 63	30.032	67.305	53.938	1.00	13.27	8
5516	OWO	WAT	V 64	32.331	68.357	33.829	1.00	13.58	8
5517	OWO	WAT	V 65	23.329	85.511	30.252	1.00	13.33	8
5518	OWO	WAT	V 66	20.246	61.387	18.981	1.00	13.61	8
5519	OWO	WAT	V 67	28.775	74.856	40.336	1.00	13.65	8
5520	OWO	WAT	V 68	32.567	68.924	43.607	1.00	13.35	8
5521	OWO	WAT	V 69	10.838	68.713	42.880	1.00	13.33	8
5522	OWO	WAT	V 70	12.859	61.518	43.606	1.00	13.90	8
5523	OWO	WAT	V 71	45.207	60.214	32.334	1.00	13.77	8
5524	OWO	WAT	V 72	27.427	66.987	53.108	1.00	13.75	8
5525	OWO	WAT	V 73	19.074	63.276	42.264	1.00	13.66	8
5526	OWO	WAT	V 74	36.934	75.592	8.270	1.00	13.53	8
5527	OWO	WAT	V 75	27.574	81.410	6.013	1.00	14.19	8
5528	OWO	WAT	V 76	30.621	83.670	31.215	1.00	14.28	8
5529	OWO	WAT	V 77	42.514	70.356	20.822	1.00	14.45	8
5530	OWO	WAT	V 78	12.529	75.168	22.815	1.00	14.52	8
5531	OWO	WAT	V 79	39.891	56.461	11.992	1.00	14.18	8
5532	OWO	WAT	V 80	30.677	68.114	68.620	1.00	14.47	8
5533	OWO	WAT	V 81	33.218	64.224	36.711	1.00	14.46	8
5534	OWO	WAT	V 82	12.035	74.811	37.533	1.00	14.63	8
5535	OWO	WAT	V 83	15.981	73.538	38.669	1.00	14.16	8
5536	OWO	WAT	V 84	10.686	81.113	35.947	1.00	14.90	8
5537	OWO	WAT	V 85	25.562	71.871	51.287	1.00	14.81	8
5538	OWO	WAT	V 86	29.447	83.564	16.644	1.00	14.86	8
5539	OWO	WAT	V 87	13.480	81.300	31.112	1.00	14.85	8
5540	OWO	WAT	V 88	5.774	80.076	41.775	1.00	14.92	8
5541	OWO	WAT	V 89	47.914	63.828	68.644	1.00	14.87	8
5542	OWO	WAT	V 90	34.743	77.662	45.101	1.00	15.02	8
5543	OWO	WAT	V 91	24.427	76.387	7.359	1.00	15.03	8
5544	OWO	WAT	V 92	-2.703	63.471	38.845	1.00	15.46	8
5545	OWO	WAT	V 93	14.681	69.092	37.625	1.00	15.14	8
5546	OWO	WAT	V 94	28.123	74.474	42.954	1.00	15.12	8
5547	OWO	WAT	V 95	23.589	80.407	29.737	1.00	15.19	8
5548	OWO	WAT	V 96	28.941	64.658	47.281	1.00	15.32	8
5549	OWO	WAT	V 97	33.848	62.923	7.441	1.00	15.09	8
5550	OWO	WAT	V 98	50.687	59.709	65.034	1.00	15.32	8
5551	OWO	WAT	V 99	29.977	80.264	14.804	1.00	15.36	8
5552	OWO	WAT	V 100	25.916	52.607	44.011	1.00	15.58	8
5553	OWO	WAT	V 101	6.765	62.176	32.083	1.00	15.56	8
5554	OWO	WAT	V 102	13.107	89.810	20.076	1.00	15.94	8
5555	OWO	WAT	V 103	11.601	78.292	36.745	1.00	15.75	8

ID								
5556	OWO	WAT	V 104	14.144	58.009	41.510	1.00	15.84	8
5557	OWO	WAT	V 105	44.010	71.623	35.990	1.00	15.89	8
5558	OWO	WAT	V 106	21.168	82.474	32.738	1.00	16.12	8
5559	OWO	WAT	V 107	28.667	48.688	46.155	1.00	16.35	8
5560	OWO	WAT	V 108	25.610	86.818	38.209	1.00	16.18	8
5561	OWO	WAT	V 109	29.070	89.992	31.492	1.00	16.47	8
5562	OWO	WAT	V 110	1.291	75.000	41.091	1.00	16.32	8
5563	OWO	WAT	V 111	34.624	56.600	71.328	1.00	16.34	8
5564	OWO	WAT	V 112	28.281	69.481	68.152	1.00	16.52	8
5565	OWO	WAT	V 113	26.135	87.038	35.304	1.00	16.58	8
5566	OWO	WAT	V 114	35.168	78.123	51.153	1.00	16.64	8
5567	OWO	WAT	V 115	19.827	81.281	46.203	1.00	16.48	8
5568	OWO	WAT	V 116	30.082	84.087	8.323	1.00	16.75	8
5569	OWO	WAT	V 117	45.164	71.238	15.992	1.00	16.58	8
5570	OWO	WAT	V 118	-2.555	86.829	29.585	1.00	16.73	8
5571	OWO	WAT	V 119	1.879	75.860	26.297	1.00	16.96	8
5572	OWO	WAT	V 120	20.960	94.415	22.713	1.00	16.87	8
5573	OWO	WAT	V 121	12.300	72.626	21.951	1.00	16.94	8
5574	OWO	WAT	V 122	21.720	86.954	37.648	1.00	17.08	8
5575	OWO	WAT	V 123	17.342	46.052	29.967	1.00	16.89	8
5576	OWO	WAT	V 124	15.847	84.337	34.562	1.00	16.74	8
5577	OWO	WAT	V 125	-3.241	60.294	45.682	1.00	16.67	8
5578	OWO	WAT	V 126	11.587	104.148	24.323	1.00	17.24	8
5579	OWO	WAT	V 127	28.501	56.852	51.441	1.00	17.40	8
5580	OWO	WAT	V 128	14.206	82.937	18.392	1.00	17.21	8
5581	OWO	WAT	V 129	41.516	56.407	63.525	1.00	17.22	8
5582	OWO	WAT	V 130	36.936	73.986	40.016	1.00	17.65	8
5583	OWO	WAT	V 131	20.790	40.115	27.752	1.00	17.53	8
5584	OWO	WAT	V 132	45.240	52.146	20.212	1.00	17.55	8
5585	OWO	WAT	V 133	41.799	49.726	28.892	1.00	17.29	8
5586	OWO	WAT	V 134	23.108	66.755	60.083	1.00	17.41	8
5587	OWO	WAT	V 135	26.863	48.805	41.435	1.00	17.67	8
5588	OWO	WAT	V 136	27.488	90.699	28.239	1.00	17.76	8
5589	OWO	WAT	V 137	0.191	69.111	49.096	1.00	17.78	8
5590	OWO	WAT	V 138	34.447	43.232	33.946	1.00	17.85	8
5591	OWO	WAT	V 139	22.589	64.044	49.415	1.00	18.18	8
5592	OWO	WAT	V 140	17.697	81.339	35.462	1.00	17.98	8
5593	OWO	WAT	V 141	2.444	96.838	37.934	1.00	18.09	8
5594	OWO	WAT	V 142	35.347	81.251	39.545	1.00	17.94	8
5595	OWO	WAT	V 143	14.511	56.297	49.652	1.00	18.35	8
5596	OWO	WAT	V 144	10.014	71.654	20.618	1.00	18.11	8
5597	OWO	WAT	V 145	47.629	63.815	18.991	1.00	18.25	8

ID								1.00		8
5598	OWO	WAT	V	146	15.832	79.962	31.635	1.00	18.52	8
5599	OWO	WAT	V	147	35.482	76.428	39.594	1.00	18.64	8
5600	OWO	WAT	V	149	16.532	67.334	48.159	1.00	18.61	8
5601	OWO	WAT	V	150	32.280	83.173	33.550	1.00	18.44	8
5602	OWO	WAT	V	151	35.037	62.885	71.628	1.00	18.62	8
5603	OWO	WAT	V	152	14.756	56.448	28.743	1.00	18.87	8
5604	OWO	WAT	V	153	51.007	64.138	32.515	1.00	18.45	8
5605	OWO	WAT	V	154	44.683	55.406	52.749	1.00	18.86	8
5606	OWO	WAT	V	155	30.803	47.286	18.413	1.00	19.00	8
5607	OWO	WAT	V	156	21.419	87.008	31.586	1.00	18.86	8
5608	OWO	WAT	V	157	27.471	54.562	13.822	1.00	18.93	8
5609	OWO	WAT	V	158	19.938	85.031	32.888	1.00	18.77	8
5610	OWO	WAT	V	160	23.159	58.927	72.012	1.00	19.01	8
5611	OWO	WAT	V	161	20.470	84.773	18.243	1.00	19.09	8
5612	OWO	WAT	V	162	15.077	66.831	13.792	1.00	18.88	8
5613	OWO	WAT	V	164	34.016	76.548	58.886	1.00	19.18	8
5614	OWO	WAT	V	165	1.791	96.077	35.490	1.00	19.39	8
5615	OWO	WAT	V	168	16.921	65.681	50.173	1.00	19.74	8
5616	OWO	WAT	V	169	36.015	49.508	63.823	1.00	19.75	8
5617	OWO	WAT	V	170	19.146	93.639	20.716	1.00	19.72	8
5618	OWO	WAT	V	171	41.081	74.262	43.114	1.00	19.99	8
5619	OWO	WAT	V	172	27.356	86.834	41.150	1.00	19.86	8
5620	OWO	WAT	V	173	32.541	61.241	6.239	1.00	19.49	8
5621	OWO	WAT	V	174	51.971	65.011	17.317	1.00	20.07	8
5622	OWO	WAT	V	175	36.754	79.161	49.332	1.00	19.96	8
5623	OWO	WAT	V	176	21.529	68.516	52.600	1.00	19.88	8
5624	OWO	WAT	V	177	52.175	64.476	48.627	1.00	19.48	8
5625	OWO	WAT	V	178	47.687	55.894	28.556	1.00	19.85	8
5626	OWO	WAT	V	180	28.309	57.733	12.302	1.00	19.86	8
5627	OWO	WAT	V	181	19.852	88.585	39.017	1.00	20.03	8
5628	OWO	WAT	V	182	48.116	65.053	11.253	1.00	19.75	8
5629	OWO	WAT	V	183	45.728	49.401	46.755	1.00	20.07	8
5630	OWO	WAT	V	184	23.090	51.403	56.173	1.00	19.76	8
5631	OWO	WAT	V	185	23.972	69.604	56.226	1.00	20.10	8
5632	OWO	WAT	V	186	49.679	65.500	55.245	1.00	19.98	8
5633	OWO	WAT	V	187	50.720	57.388	48.688	1.00	20.23	8
5634	OWO	WAT	V	188	34.857	63.152	38.228	1.00	20.32	8
5635	OWO	WAT	V	189	37.511	43.029	19.233	1.00	20.64	8
5636	OWO	WAT	V	190	50.567	67.974	10.449	1.00	20.34	8
5637	OWO	WAT	V	191	29.875	92.014	24.922	1.00	20.35	8
5638	OWO	WAT	V	192	4.055	100.811	36.567	1.00	20.35	8
5639	OWO	WAT	V	194	23.592	87.859	32.780	1.00	20.02	8

ID								1.00		8
5640	OWO	WAT	V	195	43.898	69.204	9.296	1.00	20.35	8
5641	OWO	WAT	V	196	17.503	91.553	19.944	1.00	20.36	8
5642	OWO	WAT	V	197	22.079	71.014	56.203	1.00	20.50	8
5643	OWO	WAT	V	198	24.611	90.278	32.659	1.00	20.71	8
5644	OWO	WAT	V	199	15.822	79.667	34.285	1.00	20.71	8
5645	OWO	WAT	V	200	41.507	64.519	35.769	1.00	20.47	8
5646	OWO	WAT	V	201	50.582	63.510	21.444	1.00	20.44	8
5647	OWO	WAT	V	202	-4.254	88.781	28.481	1.00	20.69	8
5648	OWO	WAT	V	203	41.289	48.120	25.349	1.00	20.50	8
5649	OWO	WAT	V	204	33.838	47.522	55.491	1.00	20.42	8
5650	OWO	WAT	V	205	28.696	42.611	30.481	1.00	21.10	8
5651	OWO	WAT	V	206	29.680	89.479	34.207	1.00	20.71	8
5652	OWO	WAT	V	207	13.375	82.989	38.528	1.00	20.79	8
5653	OWO	WAT	V	208	-0.381	94.652	35.925	1.00	20.71	8
5654	OWO	WAT	V	209	32.894	53.249	7.401	1.00	20.79	8
5655	OWO	WAT	V	210	47.202	71.009	63.961	1.00	21.03	8
5656	OWO	WAT	V	211	16.432	73.589	49.935	1.00	20.64	8
5657	OWO	WAT	V	212	36.761	42.123	23.718	1.00	20.49	8
5658	OWO	WAT	V	213	1.326	98.176	25.977	1.00	21.14	8
5659	OWO	WAT	V	215	32.369	84.143	13.694	1.00	21.27	8
5660	OWO	WAT	V	216	1.000	72.006	27.168	1.00	21.23	8
5661	OWO	WAT	V	217	21.907	50.669	62.000	1.00	21.19	8
5662	OWO	WAT	V	218	30.956	41.738	21.314	1.00	21.03	8
5663	OWO	WAT	V	219	36.121	71.463	4.684	1.00	21.46	8
5664	OWO	WAT	V	220	13.404	81.016	34.970	1.00	21.65	8
5665	OWO	WAT	V	221	22.957	65.844	62.604	1.00	21.35	8
5666	OWO	WAT	V	222	4.260	99.337	39.649	1.00	21.63	8
5667	OWO	WAT	V	223	17.535	84.263	32.555	1.00	20.98	8
5668	OWO	WAT	V	224	16.496	59.065	45.746	1.00	21.66	8
5669	OWO	WAT	V	225	16.576	52.574	34.164	1.00	21.65	8
5670	OWO	WAT	V	226	30.825	51.406	13.432	1.00	21.28	8
5671	OWO	WAT	V	227	39.177	82.848	23.121	1.00	21.63	8
5672	OWO	WAT	V	229	19.108	58.790	36.143	1.00	21.40	8
5673	OWO	WAT	V	230	19.087	75.378	50.063	1.00	21.22	8
5674	OWO	WAT	V	231	47.413	69.881	16.583	1.00	21.29	8
5675	OWO	WAT	V	232	26.686	91.040	31.058	1.00	21.61	8
5676	OWO	WAT	V	233	29.036	52.089	61.852	1.00	21.25	8
5677	OWO	WAT	V	234	36.187	79.293	46.671	1.00	21.45	8
5678	OWO	WAT	V	235	37.289	74.835	65.582	1.00	21.65	8
5679	OWO	WAT	V	236	-0.428	68.770	30.109	1.00	21.94	8
5680	OWO	WAT	V	237	50.762	57.829	44.061	1.00	22.15	8
5681	OWO	WAT	V	238	45.167	70.565	24.893	1.00	22.00	8

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5682	OWO	WAT	V	239	28.609	48.609	51.930	1.00	22.03	8
5683	OWO	WAT	V	241	2.366	94.552	38.603	1.00	21.97	8
5684	OWO	WAT	V	242	9.365	68.970	20.742	1.00	22.04	8
5685	OWO	WAT	V	243	38.583	85.642	29.447	1.00	21.90	8
5686	OWO	WAT	V	244	24.639	49.542	55.081	1.00	22.42	8
5687	OWO	WAT	V	245	18.177	95.474	18.838	1.00	22.69	8
5688	OWO	WAT	V	246	37.307	43.276	33.862	1.00	22.90	8
5689	OWO	WAT	V	247	23.478	86.290	16.425	1.00	21.85	8
5690	OWO	WAT	V	248	43.569	60.466	72.156	1.00	23.10	8
5691	OWO	WAT	V	249	23.281	40.658	26.691	1.00	22.77	8
5692	OWO	WAT	V	250	8.761	65.837	24.376	1.00	23.00	8
5693	OWO	WAT	V	251	27.215	72.803	60.509	1.00	22.60	8
5694	OWO	WAT	V	252	16.174	91.014	44.772	1.00	22.53	8
5695	OWO	WAT	V	253	3.297	98.395	35.969	1.00	23.15	8
5696	OWO	WAT	V	254	10.918	68.974	53.447	1.00	23.22	8
5697	OWO	WAT	V	255	-5.802	66.167	37.375	1.00	23.15	8
5698	OWO	WAT	V	256	8.490	91.603	44.099	1.00	22.65	8
5699	OWO	WAT	V	257	9.234	73.507	16.552	1.00	22.92	8
5700	OWO	WAT	V	258	15.730	56.257	16.556	1.00	23.08	8
5701	OWO	WAT	V	259	19.989	82.268	13.713	1.00	22.90	8
5702	OWO	WAT	V	260	27.613	49.792	60.562	1.00	23.09	8
5703	OWO	WAT	V	261	26.408	61.061	10.006	1.00	23.11	8
5704	OWO	WAT	V	262	14.277	70.834	12.781	1.00	23.32	8
5705	OWO	WAT	V	263	-3.969	72.948	34.601	1.00	22.98	8
5706	OWO	WAT	V	264	-1.028	93.829	38.475	1.00	23.17	8
5707	OWO	WAT	V	265	19.230	101.476	27.805	1.00	22.95	8
5708	OWO	WAT	V	266	17.914	54.281	55.523	1.00	23.31	8
5709	OWO	WAT	V	267	-1.163	64.024	50.895	1.00	23.41	8
5710	OWO	WAT	V	268	16.209	48.607	29.481	1.00	23.45	8
5711	OWO	WAT	V	269	35.070	63.595	75.008	1.00	23.90	8
5712	OWO	WAT	V	270	38.343	43.626	21.918	1.00	23.33	8
5713	OWO	WAT	V	271	32.197	85.153	37.181	1.00	23.49	8
5714	OWO	WAT	V	272	28.789	71.419	66.331	1.00	22.79	8
5715	OWO	WAT	V	273	41.806	79.697	18.591	1.00	23.39	8
5716	OWO	WAT	V	274	38.127	43.105	36.443	1.00	23.39	8
5717	OWO	WAT	V	275	16.104	65.021	59.364	1.00	23.44	8
5718	OWO	WAT	V	276	-6.314	69.710	39.367	1.00	24.17	8
5719	OWO	WAT	V	277	25.476	86.458	13.742	1.00	23.04	8
5720	OWO	WAT	V	278	0.680	86.883	37.755	1.00	23.27	8
5721	OWO	WAT	V	279	33.015	81.157	29.976	1.00	23.30	8
5722	OWO	WAT	V	280	1.763	88.744	19.524	1.00	23.68	8
5723	OWO	WAT	V	281	13.574	55.293	46.773	1.00	23.63	8
5724	OWO	WAT	V	282	13.480	62.142	18.325	1.00	23.48	8
5725	OWO	WAT	V	283	19.332	56.925	43.429	1.00	23.84	8
5726	OWO	WAT	V	284	7.117	53.127	42.902	1.00	23.57	8
5727	OWO	WAT	V	285	27.829	78.220	46.852	1.00	24.20	8
5728	OWO	WAT	V	286	49.295	57.105	51.124	1.00	23.77	8
5729	OWO	WAT	V	287	13.680	105.356	27.022	1.00	24.03	8
5730	OWO	WAT	V	288	17.824	106.639	32.939	1.00	24.44	8
5731	OWO	WAT	V	289	15.542	66.715	53.507	1.00	24.42	8
5732	OWO	WAT	V	290	27.050	47.324	23.573	1.00	24.19	8
5733	OWO	WAT	V	291	39.482	73.577	31.816	1.00	24.19	8
5734	OWO	WAT	V	292	10.356	77.638	16.052	1.00	23.96	8
5735	OWO	WAT	V	293	24.405	63.252	50.730	1.00	24.31	8
5736	OWO	WAT	V	294	15.639	54.776	31.091	1.00	24.22	8
5737	OWO	WAT	V	295	-5.196	74.449	41.013	1.00	24.55	8
5738	OWO	WAT	V	296	40.985	72.335	29.872	1.00	24.18	8
5739	OWO	WAT	V	297	48.449	77.051	18.858	1.00	24.84	8
5740	OWO	WAT	V	298	25.992	92.521	26.784	1.00	24.65	8
5741	OWO	WAT	V	299	45.814	49.306	39.112	1.00	24.94	8
5742	OWO	WAT	V	300	44.725	55.968	11.440	1.00	24.73	8
5743	OWO	WAT	V	301	20.058	84.809	46.357	1.00	24.58	8
5744	OWO	WAT	V	302	18.079	50.984	17.334	1.00	24.81	8
5745	OWO	WAT	V	303	17.020	65.666	56.879	1.00	24.66	8
5746	OWO	WAT	V	304	44.682	72.661	22.319	1.00	24.37	8
5747	OWO	WAT	V	305	0.091	75.457	45.401	1.00	24.90	8
5748	OWO	WAT	V	306	50.222	58.393	53.380	1.00	25.29	8
5749	OWO	WAT	V	307	44.639	52.674	53.486	1.00	24.60	8
5750	OWO	WAT	V	308	49.725	64.827	52.493	1.00	24.82	8
5751	OWO	WAT	V	309	39.542	61.118	7.698	1.00	25.15	8
5752	OWO	WAT	V	310	41.190	79.148	23.073	1.00	24.95	8
5753	OWO	WAT	V	311	1.598	80.899	39.623	1.00	24.92	8
5754	OWO	WAT	V	312	25.053	45.139	19.817	1.00	25.20	8
5755	OWO	WAT	V	313	16.135	84.191	47.663	1.00	25.29	8
5756	OWO	WAT	V	314	11.381	71.857	49.110	1.00	25.43	8
5757	OWO	WAT	V	315	-4.512	57.278	42.326	1.00	25.09	8
5758	OWO	WAT	V	316	3.805	93.061	22.790	1.00	25.16	8
5759	OWO	WAT	V	317	34.832	82.782	37.221	1.00	25.31	8
5760	OWO	WAT	V	318	3.711	95.706	23.179	1.00	25.69	8
5761	OWO	WAT	V	319	20.209	70.499	54.469	1.00	25.01	8
5762	OWO	WAT	V	320	11.697	68.339	56.789	1.00	25.40	8
5763	OWO	WAT	V	321	7.254	91.019	18.448	1.00	25.81	8
5764	OWO	WAT	V	322	42.661	49.979	21.111	1.00	25.51	8
5765	OWO	WAT	V	323	17.078	70.892	50.293	1.00	25.51	8

ID	Label					
5766	OWO WAT V 324	49.200	60.915	31.600	1.00	25.53	8
5767	OWO WAT V 325	31.661	71.214	4.552	1.00	25.64	8
5768	OWO WAT V 326	-2.062	96.310	27.990	1.00	25.87	8
5769	OWO WAT V 327	39.125	80.230	56.164	1.00	26.25	8
5770	OWO WAT V 328	48.433	70.294	57.525	1.00	25.87	8
5771	OWO WAT V 329	-2.772	64.990	34.941	1.00	25.71	8
5772	OWO WAT V 330	35.352	40.001	22.353	1.00	25.96	8
5773	OWO WAT V 331	34.557	81.627	46.792	1.00	26.23	8
5774	OWO WAT V 332	23.250	92.581	31.676	1.00	26.20	8
5775	OWO WAT V 333	23.167	53.251	13.715	1.00	25.74	8
5776	OWO WAT V 334	20.707	52.356	39.202	1.00	25.65	8
5777	OWO WAT V 335	4.870	93.727	41.264	1.00	26.54	8
5778	OWO WAT V 336	0.053	71.401	29.885	1.00	25.67	8
5779	OWO WAT V 337	20.015	56.517	72.324	1.00	25.86	8
5780	OWO WAT V 338	13.826	74.132	49.507	1.00	26.36	8
5781	OWO WAT V 339	44.958	71.632	28.685	1.00	26.80	8
5782	OWO WAT V 340	-0.093	73.784	33.119	1.00	26.07	8
5783	OWO WAT V 341	49.630	56.579	40.106	1.00	26.64	8
5784	OWO WAT V 342	15.628	79.186	45.524	1.00	26.42	8
5785	OWO WAT V 343	31.176	91.867	30.769	1.00	26.10	8
5786	OWO WAT V 344	15.626	67.643	55.543	1.00	26.46	8
5787	OWO WAT V 345	21.398	95.678	38.839	1.00	26.79	8
5788	OWO WAT V 346	41.099	41.982	39.915	1.00	26.66	8
5789	OWO WAT V 347	22.442	79.888	2.661	1.00	26.18	8
5790	OWO WAT V 348	44.448	72.532	39.099	1.00	27.08	8
5791	OWO WAT V 349	40.265	82.719	10.175	1.00	26.27	8
5792	OWO WAT V 350	40.934	43.657	30.662	1.00	27.04	8
5793	OWO WAT V 351	-1.666	97.213	32.163	1.00	27.31	8
5794	OWO WAT V 352	-8.111	67.221	45.293	1.00	27.12	8
5795	OWO WAT V 353	16.355	56.614	42.471	1.00	27.10	8
5796	OWO WAT V 354	11.346	65.531	19.934	1.00	27.40	8
5797	OWO WAT V 355	11.189	105.275	21.949	1.00	27.29	8
5798	OWO WAT V 356	23.545	83.502	8.615	1.00	27.06	8
5799	OWO WAT V 357	22.122	49.913	14.384	1.00	27.06	8
5800	OWO WAT V 358	6.833	52.668	34.000	1.00	27.79	8
5801	OWO WAT V 359	30.479	47.598	56.666	1.00	27.66	8
5802	OWO WAT V 360	33.166	42.826	19.802	1.00	27.33	8
5803	OWO WAT V 361	24.029	55.851	13.947	1.00	28.08	8
5804	OWO WAT V 362	39.488	85.709	20.632	1.00	27.01	8
5805	OWO WAT V 363	2.130	49.916	43.049	1.00	27.26	8
5806	OWO WAT V 364	35.616	41.373	27.626	1.00	27.13	8
5807	OWO WAT V 365	50.664	58.560	41.450	1.00	26.59	8

ID	Label					
5808	OWO WAT V 366	30.943	49.068	15.778	1.00	26.98	8
5809	OWO WAT V 367	43.193	44.166	34.878	1.00	26.97	8
5810	OWO WAT V 368	6.994	74.430	21.876	1.00	27.26	8
5811	OWO WAT V 369	34.427	88.111	38.939	1.00	27.30	8
5812	OWO WAT V 370	1.636	93.582	40.791	1.00	27.54	8
5813	OWO WAT V 371	5.971	98.241	21.504	1.00	27.42	8
5814	OWO WAT V 372	29.223	75.038	60.445	1.00	26.93	8
5815	OWO WAT V 373	31.316	53.650	9.481	1.00	27.85	8
5816	OWO WAT V 374	43.939	56.548	55.050	1.00	27.63	8
5817	OWO WAT V 375	46.559	74.427	54.610	1.00	27.46	8
5818	OWO WAT V 376	26.961	70.400	70.205	1.00	28.10	8
5819	OWO WAT V 377	48.989	57.080	35.445	1.00	28.37	8
5820	OWO WAT V 378	14.950	47.050	27.899	1.00	27.90	8
5821	OWO WAT V 379	46.760	46.713	39.616	1.00	27.67	8
5822	OWO WAT V 380	-1.380	79.896	38.771	1.00	28.02	8
5823	OWO WAT V 381	12.690	81.625	15.826	1.00	28.03	8
5824	OWO WAT V 382	10.617	104.039	32.646	1.00	28.19	8
5825	OWO WAT V 383	13.859	80.594	13.625	1.00	27.99	8
5826	OWO WAT V 384	7.322	72.231	20.822	1.00	28.35	8
5827	OWO WAT V 385	29.284	46.632	44.658	1.00	28.26	8
5828	OWO WAT V 386	18.064	50.559	24.862	1.00	28.35	8
5829	OWO WAT V 387	35.054	45.412	46.835	1.00	28.58	8
5830	OWO WAT V 388	22.478	83.558	13.277	1.00	28.32	8
5831	OWO WAT V 389	10.928	67.510	14.236	1.00	28.65	8
5832	OWO WAT V 390	33.397	82.246	49.220	1.00	27.37	8
5833	OWO WAT V 391	23.434	88.556	18.508	1.00	28.95	8
5834	OWO WAT V 392	29.832	42.212	39.251	1.00	28.21	8
5835	OWO WAT V 393	15.076	70.102	51.758	1.00	27.91	8
5836	OWO WAT V 394	35.566	48.453	66.263	1.00	28.16	8
5837	OWO WAT V 395	34.691	46.594	49.301	1.00	28.85	8
5838	OWO WAT V 396	39.702	74.440	36.238	1.00	28.77	8
5839	OWO WAT V 397	50.176	56.086	14.851	1.00	28.75	8
5840	OWO WAT V 398	14.115	57.834	59.211	1.00	27.83	8
5841	OWO WAT V 399	38.189	49.702	11.383	1.00	29.29	8
5842	OWO WAT V 400	3.241	100.941	39.033	1.00	28.47	8
5843	OWO WAT V 401	44.298	73.510	63.099	1.00	28.91	8
5844	OWO WAT V 402	7.061	75.135	47.544	1.00	28.59	8
5845	OWO WAT V 403	5.477	86.910	42.452	1.00	29.56	8
5846	OWO WAT V 404	25.564	72.515	56.246	1.00	29.09	8
5847	OWO WAT V 405	25.075	41.160	32.189	1.00	29.20	8
5848	OWO WAT V 406	2.074	56.387	58.854	1.00	28.86	8
5849	OWO WAT V 407	47.163	58.038	33.355	1.00	28.76	8

No.	Label						
5850	OWO WAT V 408	27.438	44.407	21.116	1.00	28.21	8
5851	OWO WAT V 409	32.690	89.765	34.445	1.00	29.13	8
5852	OWO WAT V 410	8.726	56.984	33.875	1.00	28.81	8
5853	OWO WAT V 411	-2.765	74.847	45.763	1.00	28.58	8
5854	OWO WAT V 412	9.880	76.028	48.593	1.00	29.43	8
5855	OWO WAT V 413	18.587	55.405	41.191	1.00	28.66	8
5856	OWO WAT V 414	0.831	95.595	23.928	1.00	28.90	8
5857	OWO WAT V 415	18.167	58.042	65.360	1.00	28.57	8
5858	OWO WAT V 416	42.814	47.661	27.620	1.00	29.34	8
5859	OWO WAT V 417	19.226	89.316	19.020	1.00	29.13	8
5860	OWO WAT V 418	17.933	82.283	11.952	1.00	28.48	8
5861	OWO WAT V 419	44.723	79.313	13.493	1.00	30.12	8
5862	OWO WAT V 420	34.399	78.316	0.819	1.00	29.81	8
5863	OWO WAT V 421	28.282	72.460	57.888	1.00	29.12	8
5864	OWO WAT V 422	50.448	63.547	18.793	1.00	29.31	8
5865	OWO WAT V 423	43.033	55.323	65.913	1.00	29.32	8
5866	OWO WAT V 424	45.865	51.168	51.498	1.00	29.29	8
5867	OWO WAT V 425	12.844	59.751	19.386	1.00	29.45	8
5868	OWO WAT V 426	-3.707	82.697	22.959	1.00	29.58	8
5869	OWO WAT V 427	28.989	45.171	18.941	1.00	29.97	8
5870	OWO WAT V 428	14.689	65.149	11.472	1.00	29.76	8
5871	OWO WAT V 429	49.962	74.061	53.225	1.00	29.59	8
5872	OWO WAT V 430	26.674	46.428	17.647	1.00	29.59	8
5873	OWO WAT V 431	32.017	64.281	73.846	1.00	29.67	8
5874	OWO WAT V 432	-1.685	89.735	36.396	1.00	29.42	8
5875	OWO WAT V 433	35.326	90.671	32.858	1.00	30.08	8
5876	OWO WAT V 434	-5.217	86.519	22.737	1.00	30.64	8
5877	OWO WAT V 435	22.655	40.619	31.468	1.00	30.18	8
5878	OWO WAT V 436	9.970	66.686	21.940	1.00	30.16	8
5879	OWO WAT V 437	22.156	103.267	41.579	1.00	28.97	8
5880	OWO WAT V 438	40.919	45.553	23.996	1.00	30.25	8
5881	OWO WAT V 439	24.378	42.994	18.498	1.00	29.87	8
5882	OWO WAT V 440	27.456	90.485	38.760	1.00	30.24	8
5883	OWO WAT V 441	41.628	78.372	57.406	1.00	30.51	8
5884	OWO WAT V 442	10.162	67.365	58.879	1.00	29.98	8
5885	OWO WAT V 443	7.899	70.258	55.642	1.00	30.04	8
5886	OWO WAT V 444	7.439	62.721	5.170	1.00	30.62	8
5887	OWO WAT V 445	30.182	60.296	16.010	1.00	30.22	8
5888	OWO WAT V 446	36.835	87.065	55.789	1.00	30.39	8
5889	OWO WAT V 447	48.154	58.076	38.708	1.00	30.45	8
5890	OWO WAT V 448	9.079	102.838	5.903	1.00	30.35	8
5891	OWO WAT V 449	31.488	67.009		1.00		8

No.	Label						
5892	OWO WAT V 450	8.486	84.645	16.251	1.00	30.46	8
5893	OWO WAT V 451	17.712	70.200	52.856	1.00	31.30	8
5894	OWO WAT V 452	45.187	54.640	9.218	1.00	30.47	8
5895	OWO WAT V 453	23.220	65.098	2.001	1.00	30.15	8
5896	OWO WAT V 454	43.353	48.993	23.792	1.00	30.21	8
5897	OWO WAT V 455	0.241	76.716	42.809	1.00	30.43	8
5898	OWO WAT V 456	38.954	90.362	25.885	1.00	31.19	8
5899	OWO WAT V 457	8.998	51.618	42.193	1.00	30.65	8
5900	OWO WAT V 458	47.484	70.509	20.484	1.00	30.75	8
5901	OWO WAT V 459	26.632	83.732	7.006	1.00	30.86	8
5902	OWO WAT V 460	27.887	84.031	46.702	1.00	31.16	8
5903	OWO WAT V 461	10.456	52.659	39.931	1.00	31.37	8
5904	OWO WAT V 462	25.474	53.247	11.551	1.00	31.54	8
5905	OWO WAT V 463	21.666	48.674	41.178	1.00	31.71	8
5906	OWO WAT V 464	51.799	63.643	37.234	1.00	30.69	8
5907	OWO WAT V 465	17.686	48.668	36.134	1.00	31.38	8
5908	OWO WAT V 466	47.081	52.951	49.717	1.00	30.51	8
5909	OWO WAT V 467	15.593	93.203	42.768	1.00	30.90	8
5910	OWO WAT V 468	36.480	80.071	53.771	1.00	30.56	8
5911	OWO WAT V 469	42.137	47.781	14.882	1.00	31.85	8
5912	OWO WAT V 470	22.333	61.351	70.096	1.00	31.36	8
5913	OWO WAT V 471	40.228	72.527	72.095	1.00	31.17	8
5914	OWO WAT V 472	1.666	74.527	48.132	1.00	31.28	8
5915	OWO WAT V 473	14.531	66.936	59.718	1.00	31.26	8
5916	OWO WAT V 474	21.000	68.421	1.879	1.00	31.59	8
5917	OWO WAT V 475	37.163	82.170	41.385	1.00	31.32	8
5918	OWO WAT V 476	26.307	62.094	4.445	1.00	31.75	8
5919	OWO WAT V 477	51.579	61.409	40.641	1.00	31.25	8
5920	OWO WAT V 478	41.761	64.533	7.578	1.00	31.59	8
5921	OWO WAT V 479	-7.353	65.632	35.219	1.00	31.13	8
5922	OWO WAT V 480	16.514	82.720	36.237	1.00	31.25	8
5923	OWO WAT V 481	23.770	75.891	54.586	1.00	31.53	8
5924	OWO WAT V 482	50.377	68.458	56.245	1.00	31.72	8
5925	OWO WAT V 483	15.979	105.053	38.501	1.00	30.98	8
5926	OWO WAT V 484	27.194	93.867	31.829	1.00	31.85	8
5927	OWO WAT V 485	39.487	68.664	5.438	1.00	31.02	8
5928	OWO WAT V 486	30.446	58.438	75.573	1.00	31.49	8
5929	OWO WAT V 487	-2.258	85.349	34.418	1.00	31.74	8
5930	OWO WAT V 488	46.873	56.769	65.770	1.00	31.57	8
5931	OWO WAT V 489	18.407	47.617	19.229	1.00	31.61	8
5932	OWO WAT V 490	2.500	78.686	43.392	1.00	31.62	8
5933	OWO WAT V 491	7.727	97.913	42.497	1.00	32.36	8

ID	Name		val1	val2	val3		val4	
5934	OWO WAT V	492	-5.430	71.658	30.991	1.00	32.45	8
5935	OWO WAT V	493	47.547	72.958	21.222	1.00	31.57	8
5936	OWO WAT V	494	9.955	91.961	24.393	1.00	32.43	8
5937	OWO WAT V	495	12.996	52.420	46.232	1.00	31.76	8
5938	OWO WAT V	496	7.952	65.594	58.841	1.00	32.02	8
5939	OWO WAT V	497	37.204	88.789	33.415	1.00	31.57	8
5940	OWO WAT V	498	36.857	41.938	29.865	1.00	32.38	8
5941	OWO WAT V	499	7.220	51.844	54.727	1.00	32.32	8
5942	OWO WAT V	500	16.110	76.377	42.135	1.00	32.31	8
5943	OWO WAT V	501	24.511	47.736	48.051	1.00	32.37	8
5944	OWO WAT V	502	22.783	46.357	7.569	1.00	32.64	8
5945	OWO WAT V	503	27.138	60.897	36.850	1.00	32.20	8
5946	OWO WAT V	504	47.227	50.290	23.991	1.00	32.13	8
5947	OWO WAT V	505	6.733	67.493	56.941	1.00	32.68	8
5948	OWO WAT V	506	16.514	70.176	8.514	1.00	32.03	8
5949	OWO WAT V	507	43.175	79.662	22.899	1.00	31.76	8
5950	OWO WAT V	508	-3.757	80.394	45.262	1.00	32.62	8
5951	OWO WAT V	509	10.932	85.936	41.755	1.00	32.65	8
5952	OWO WAT V	510	3.455	97.650	69.455	1.00	32.31	8
5953	OWO WAT V	511	33.419	50.551	23.550	1.00	32.14	8
5954	OWO WAT V	512	4.069	99.962	41.690	1.00	33.21	8
5955	OWO WAT V	513	-7.206	65.717	61.406	1.00	33.20	8
5956	OWO WAT V	514	53.785	60.262	17.346	1.00	33.37	8
5957	OWO WAT V	515	16.599	61.406	30.752	1.00	32.99	8
5958	OWO WAT V	516	47.349	54.074	42.409	1.00	32.14	8
5959	OWO WAT V	517	7.038	85.163	27.538	1.00	32.78	8
5960	OWO WAT V	518	16.879	50.503	22.385	1.00	32.69	8
5961	OWO WAT V	519	30.838	91.776	32.218	1.00	33.06	8
5962	OWO WAT V	520	33.882	41.296	23.358	1.00	34.59	8
5963	OWO WAT V	521	53.502	60.512	34.316	1.00	32.27	8
5964	OWO WAT V	522	41.841	73.003	5.371	1.00	33.63	8
5965	OWO WAT V	523	18.807	69.394	27.670	1.00	32.65	8
5966	OWO WAT V	524	0.567	102.365	64.451	1.00	32.87	8
5967	OWO WAT V	525	28.899	48.822	21.339	1.00	32.62	8
5968	OWO WAT V	526	42.766	46.627	60.034	1.00	33.04	8
5969	OWO WAT V	527	18.159	66.472	59.427	1.00	33.02	8
5970	OWO WAT V	528	22.385	49.518	35.438	1.00	33.48	8
5971	OWO WAT V	529	14.542	53.899	18.821	1.00	32.40	8
5972	OWO WAT V	530	0.923	81.260	25.073	1.00	33.19	8
5973	OWO WAT V	531	35.980	91.602	71.313	1.00	32.89	8
5974	OWO WAT V	532	40.667	67.024	60.713	1.00	32.79	8
5975	OWO WAT V	533	8.626	61.294		1.00	32.69	8

ID	Name		val1	val2	val3		val4	
5976	OWO WAT V	534	35.615	48.266	53.410	1.00	33.74	8
5977	OWO WAT V	535	50.581	54.379	41.752	1.00	32.75	8
5978	OWO WAT V	536	34.738	50.213	7.372	1.00	33.68	8
5979	OWO WAT V	537	42.324	79.266	54.340	1.00	32.72	8
5980	OWO WAT V	538	16.132	78.441	8.864	1.00	33.64	8
5981	OWO WAT V	539	20.256	45.911	20.702	1.00	33.89	8
5982	OWO WAT V	540	37.253	43.882	46.170	1.00	33.73	8
5983	OWO WAT V	541	31.158	88.960	12.651	1.00	34.15	8
5984	OWO WAT V	542	2.885	65.252	30.182	1.00	33.66	8
5985	OWO WAT V	543	24.093	72.008	54.068	1.00	33.53	8
5986	OWO WAT V	544	-4.660	58.902	52.534	1.00	34.13	8
5987	OWO WAT V	545	20.523	52.571	71.335	1.00	34.29	8
5988	OWO WAT V	546	50.389	70.373	43.120	1.00	34.05	8
5989	OWO WAT V	547	-1.784	74.917	34.717	1.00	34.27	8
5990	OWO WAT V	548	25.051	100.468	31.517	1.00	33.73	8
5991	OWO WAT V	549	21.989	83.148	46.194	1.00	34.71	8
5992	OWO WAT V	550	47.521	62.364	11.718	1.00	34.05	8
5993	OWO WAT V	551	52.236	59.305	62.786	1.00	33.21	8
5994	OWO WAT V	552	40.232	80.510	16.448	1.00	34.62	8
5995	OWO WAT V	553	46.253	56.949	70.930	1.00	33.53	8
5996	OWO WAT V	554	47.895	53.053	37.454	1.00	33.41	8
5997	OWO WAT V	555	13.358	71.030	49.662	1.00	34.43	8
5998	OWO WAT V	556	-0.137	73.990	26.087	1.00	34.05	8
5999	OWO WAT V	557	43.973	74.382	6.399	1.00	34.44	8
6000	OWO WAT V	558	35.593	60.367	77.022	1.00	33.63	8
6001	OWO WAT V	559	6.112	83.592	41.192	1.00	34.97	8
6002	OWO WAT V	560	38.614	41.425	25.492	1.00	34.42	8
6003	OWO WAT V	561	34.074	88.144	12.960	1.00	34.41	8
6004	OWO WAT V	562	40.114	63.213	5.786	1.00	34.56	8
6005	OWO WAT V	563	-0.202	64.215	60.019	1.00	34.60	8
6006	OWO WAT V	564	4.614	92.264	20.329	1.00	33.82	8
6007	OWO WAT V	565	15.212	50.738	32.687	1.00	34.54	8
6008	OWO WAT V	566	13.018	70.753	53.378	1.00	35.00	8
6009	OWO WAT V	567	37.836	57.700	75.968	1.00	33.91	8
6010	OWO WAT V	568	18.054	71.354	7.047	1.00	34.28	8
6011	OWO WAT V	569	20.435	57.091	12.782	1.00	34.06	8
6012	OWO WAT V	570	-1.113	72.836	30.802	1.00	34.18	8
6013	OWO WAT V	571	45.394	47.985	15.576	1.00	34.25	8
6014	OWO WAT V	572	18.083	105.524	25.739	1.00	35.26	8
6015	OWO WAT V	573	42.363	51.035	61.868	1.00	34.64	8
6016	OWO WAT V	574	3.332	65.717	27.045	1.00	35.54	8
6017	OWO WAT V	575	0.099	70.672	51.291	1.00	34.72	8

ID	Label						
6018	OWO WAT V 576	0.045	61.496	31.457	1.00	35.51	8
6019	OWO WAT V 577	46.395	53.881	61.040	1.00	34.53	8
6020	OWO WAT V 578	52.136	68.150	59.956	1.00	35.10	8
6021	OWO WAT V 579	19.009	41.546	26.130	1.00	33.65	8
6022	OWO WAT V 580	34.720	74.712	71.942	1.00	34.54	8
6023	OWO WAT V 581	37.259	73.783	73.743	1.00	35.82	8
6024	OWO WAT V 582	27.833	55.518	10.945	1.00	34.54	8
6025	OWO WAT V 583	46.309	60.566	60.060	1.00	35.04	8
6026	OWO WAT V 584	10.952	63.329	18.773	1.00	34.23	8
6027	OWO WAT V 585	-1.056	54.481	35.620	1.00	34.65	8
6028	OWO WAT V 586	-4.178	63.600	36.602	1.00	35.48	8
6029	OWO WAT V 587	4.690	96.750	20.660	1.00	34.69	8
6030	OWO WAT V 588	8.484	73.376	50.572	1.00	35.07	8
6031	OWO WAT V 589	44.017	50.506	11.023	1.00	35.68	8
6032	OWO WAT V 590	3.744	50.549	36.911	1.00	35.69	8
6033	OWO WAT V 591	49.672	55.539	26.394	1.00	35.74	8
6034	OWO WAT V 592	27.932	93.666	25.261	1.00	35.64	8
6035	OWO WAT V 593	22.451	73.996	54.266	1.00	35.51	8
6036	OWO WAT V 594	33.048	48.697	67.743	1.00	35.66	8
6037	OWO WAT V 595	12.820	101.046	44.339	1.00	36.16	8
6038	OWO WAT V 596	51.198	69.586	53.643	1.00	35.14	8
6039	OWO WAT V 597	29.662	79.574	54.942	1.00	35.66	8
6040	OWO WAT V 598	32.247	80.128	52.116	1.00	36.54	8
6041	OWO WAT V 599	6.142	51.551	37.685	1.00	35.34	8
6042	OWO WAT V 600	5.622	70.631	22.637	1.00	35.66	8
6043	OWO WAT V 601	22.365	48.081	48.242	1.00	36.08	8
6044	OWO WAT V 602	8.288	81.332	15.641	1.00	35.59	8
6045	OWO WAT V 603	10.941	74.805	15.117	1.00	35.60	8
6046	OWO WAT V 604	35.796	73.953	1.142	1.00	35.09	8
6047	OWO WAT V 605	38.777	85.286	9.309	1.00	36.46	8
6048	OWO WAT V 606	20.560	52.963	75.539	1.00	37.03	8
6049	OWO WAT V 607	-3.377	66.230	32.749	1.00	35.54	8
6050	OWO WAT V 608	40.087	44.107	14.031	1.00	36.33	8
6051	OWO WAT V 609	-1.635	86.958	36.864	1.00	35.23	8
6052	OWO WAT V 610	48.364	69.799	29.088	1.00	35.08	8
6053	OWO WAT V 611	18.383	50.939	35.706	1.00	36.23	8
6054	OWO WAT V 612	18.334	58.135	13.735	1.00	37.22	8
6055	OWO WAT V 613	16.341	55.369	50.997	1.00	35.78	8
6056	OWO WAT V 614	28.580	52.384	12.139	1.00	36.07	8
6057	OWO WAT V 615	23.108	103.615	38.927	1.00	36.81	8
6058	OWO WAT V 616	12.325	79.803	46.301	1.00	35.61	8
6059	OWO WAT V 617	21.280	67.375	64.153	1.00	36.55	8
6060	OWO WAT V 618	40.907	80.214	52.034	1.00	37.09	8
6061	OWO WAT V 619	47.006	72.793	58.223	1.00	36.47	8
6062	OWO WAT V 620	20.221	90.473	42.499	1.00	36.27	8
6063	OWO WAT V 621	4.100	73.674	21.447	1.00	36.94	8
6064	OWO WAT V 622	8.503	104.300	35.973	1.00	36.81	8
6065	OWO WAT V 623	30.691	78.729	57.692	1.00	35.94	8
6066	OWO WAT V 624	-3.070	94.866	30.332	1.00	36.58	8
6067	OWO WAT V 625	46.455	79.736	18.637	1.00	35.97	8
6068	OWO WAT V 626	46.537	55.701	59.420	1.00	37.10	8
6069	OWO WAT V 627	43.765	47.227	50.091	1.00	37.02	8
6070	OWO WAT V 628	-0.451	52.302	37.499	1.00	36.00	8
6071	OWO WAT V 629	47.301	50.088	41.392	1.00	37.12	8
6072	OWO WAT V 630	37.956	80.815	13.060	1.00	37.94	8
6073	OWO WAT V 631	17.138	62.973	7.515	1.00	37.42	8
6074	OWO WAT V 632	13.837	57.372	18.395	1.00	36.82	8
6075	OWO WAT V 633	41.756	55.821	70.896	1.00	36.92	8
6076	OWO WAT V 634	25.077	93.998	28.706	1.00	36.97	8
6077	OWO WAT V 635	47.608	54.704	64.130	1.00	37.13	8
6078	OWO WAT V 636	6.911	105.878	25.141	1.00	36.23	8
6079	OWO WAT V 637	42.682	67.557	69.041	1.00	36.92	8
6080	OWO WAT V 638	15.363	88.594	48.034	1.00	37.42	8
6081	OWO WAT V 639	33.890	42.321	17.015	1.00	37.23	8
6082	OWO WAT V 640	9.736	84.033	43.853	1.00	36.17	8
6083	OWO WAT V 641	10.886	51.035	46.103	1.00	37.66	8
6084	OWO WAT V 642	27.929	72.978	70.610	1.00	37.35	8
6085	OWO WAT V 643	1.233	69.664	25.786	1.00	37.56	8
6086	OWO WAT V 644	15.151	107.231	33.216	1.00	36.50	8
6087	OWO WAT V 645	25.823	98.361	30.305	1.00	36.65	8
6088	OWO WAT V 646	34.857	84.394	12.899	1.00	36.71	8
6089	OWO WAT V 647	20.425	52.936	41.877	1.00	36.99	8
6090	OWO WAT V 648	26.575	49.377	65.800	1.00	37.37	8
6091	OWO WAT V 649	31.193	44.914	15.071	1.00	36.96	8
6092	OWO WAT V 650	10.578	49.772	48.688	1.00	36.01	8
6093	OWO WAT V 651	18.128	51.717	38.409	1.00	37.57	8
6094	OWO WAT V 652	45.424	54.692	56.844	1.00	37.37	8
6095	OWO WAT V 653	28.011	90.646	18.192	1.00	37.91	8
6096	OWO WAT V 654	47.536	55.043	51.941	1.00	36.73	8
6097	OWO WAT V 655	24.853	98.451	32.915	1.00	37.70	8
6098	OWO WAT V 656	27.088	98.198	26.870	1.00	38.56	8
6099	OWO WAT V 657	33.663	49.796	11.597	1.00	37.19	8
6100	OWO WAT V 658	16.309	55.242	45.035	1.00	37.83	8
6101	OWO WAT V 659	20.804	43.210	25.195	1.00	37.72	8

6102	OWO WAT V 660	23.932	65.246	70.893	1.00	37.09	8
6103	OWO WAT V 661	0.664	53.286	53.271	1.00	38.36	8
6104	OWO WAT V 662	40.187	84.151	18.538	1.00	38.05	8
6105	OWO WAT V 663	44.129	77.826	7.095	1.00	38.36	8
6106	OWO WAT V 664	5.654	103.176	36.651	1.00	38.12	8
6107	OWO WAT V 665	-2.551	55.092	37.581	1.00	39.63	8
6108	OWO WAT V 666	48.102	56.702	13.267	1.00	37.64	8
6109	OWO WAT V 667	27.824	90.475	15.555	1.00	36.41	8
6110	OWO WAT V 668	38.699	49.812	63.112	1.00	37.80	8
6111	OWO WAT V 669	-2.199	54.401	53.588	1.00	39.21	8
6112	OWO WAT V 670	19.006	53.240	47.228	1.00	38.93	8
6113	OWO WAT V 671	46.897	48.817	19.829	1.00	39.11	8
6114	OWO WAT V 672	42.301	48.935	12.664	1.00	37.43	8
6115	OWO WAT V 673	16.098	53.972	53.322	1.00	38.55	8
6116	OWO WAT V 674	16.992	66.360	61.965	1.00	40.47	8
6117	OWO WAT V 675	17.396	84.259	14.592	1.00	37.45	8
6118	OWO WAT V 676	48.023	70.300	38.414	1.00	38.28	8
6119	OWO WAT V 677	21.061	60.355	72.008	1.00	38.68	8
6120	OWO WAT V 678	25.412	45.777	15.102	1.00	40.05	8
6121	OWO WAT V 679	43.705	73.272	1.624	1.00	38.78	8
6122	OWO WAT V 680	17.379	67.334	7.173	1.00	38.88	8
6123	OWO WAT V 681	19.500	72.583	5.113	1.00	41.12	8
6124	OWO WAT V 682	27.135	53.269	75.096	1.00	38.74	8
6125	OWO WAT V 683	43.871	76.670	2.397	1.00	38.42	8
6126	OWO WAT V 684	32.690	47.277	60.342	1.00	38.23	8
6127	OWO WAT V 685	38.889	50.643	67.387	1.00	38.05	8
6128	OWO WAT V 686	15.848	75.816	7.403	1.00	37.86	8
6129	OWO WAT V 687	25.036	96.149	30.003	1.00	38.02	8
6130	OWO WAT V 688	44.120	49.147	54.661	1.00	39.21	8
6131	OWO WAT V 689	31.793	51.208	10.937	1.00	39.53	8
6132	OWO WAT V 690	9.476	96.492	43.783	1.00	37.31	8
6133	OWO WAT V 691	8.821	88.129	44.638	1.00	39.95	8
6134	OWO WAT V 692	8.003	80.636	43.644	1.00	37.70	8
6135	OWO WAT V 693	14.834	83.314	13.883	1.00	39.80	8
6136	OWO WAT V 694	21.370	44.389	41.265	1.00	39.75	8
6137	OWO WAT V 695	25.208	85.122	7.995	1.00	38.90	8
6138	OWO WAT V 696	34.431	64.498	5.882	1.00	37.43	8
6139	OWO WAT V 697	46.857	48.623	44.260	1.00	38.67	8
6140	OWO WAT V 698	48.394	49.074	31.806	1.00	38.40	8
6141	OWO WAT V 699	6.330	71.875	18.323	1.00	39.24	8
6142	OWO WAT V 700	3.567	68.493	56.513	1.00	39.03	8
6143	OWO WAT V 701	19.531	67.580	61.273	1.00	38.68	8
6144	OWO WAT V 702	36.514	82.654	34.129	1.00	38.10	8
6145	OWO WAT V 703	43.559	50.161	59.718	1.00	39.15	8
6146	OWO WAT V 704	41.881	70.064	7.497	1.00	39.49	8
6147	OWO WAT V 705	12.039	106.892	30.442	1.00	39.69	8
6148	OWO WAT V 706	25.673	81.581	52.252	1.00	39.17	8
6149	OWO WAT V 707	46.855	74.915	26.181	1.00	40.35	8
6150	OWO WAT V 708	21.077	93.052	39.231	1.00	39.44	8
6151	OWO WAT V 709	28.621	62.328	5.470	1.00	38.56	8
6152	OWO WAT V 710	-3.607	57.742	54.506	1.00	38.82	8
6153	OWO WAT V 711	25.913	41.714	29.843	1.00	38.94	8
6154	OWO WAT V 712	21.385	95.763	42.160	1.00	39.85	8
6155	OWO WAT V 713	39.666	46.565	56.093	1.00	40.16	8
6156	OWO WAT V 714	33.043	47.398	12.747	1.00	38.56	8
6157	OWO WAT V 715	3.335	48.480	47.523	1.00	40.24	8
6158	OWO WAT V 716	-5.954	59.355	43.289	1.00	39.30	8
6159	OWO WAT V 717	-2.633	70.380	52.100	1.00	39.49	8
6160	OWO WAT V 718	14.943	51.881	28.927	1.00	41.17	8
6161	OWO WAT V 719	30.665	55.789	76.270	1.00	36.29	8
6162	OWO WAT V 720	38.153	41.243	44.336	1.00	39.61	8
6163	OWO WAT V 721	42.857	72.924	4.760	1.00	39.33	8
6164	OWO WAT V 722	42.983	45.506	30.086	1.00	39.35	8
6165	OWO WAT V 723	8.541	68.671	15.477	1.00	40.10	8
6166	OWO WAT V 724	3.357	68.745	24.682	1.00	39.21	8
6167	OWO WAT V 725	32.096	74.899	2.155	1.00	38.68	8
6168	OWO WAT V 726	-3.316	99.677	31.930	1.00	40.29	8
6169	OWO WAT V 727	33.290	44.355	45.552	1.00	39.94	8
6170	OWO WAT V 728	50.677	69.314	25.213	1.00	38.89	8
6171	OWO WAT V 729	35.441	42.395	37.048	1.00	39.29	8
6172	OWO WAT V 730	8.363	53.922	35.516	1.00	38.32	8
6173	OWO WAT V 731	-1.713	91.588	39.071	1.00	40.09	8
6174	OWO WAT V 732	22.068	82.950	49.452	1.00	38.86	8
6175	OWO WAT V 733	-1.449	89.011	21.652	1.00	39.99	8
6176	OWO WAT V 734	37.354	59.035	5.289	1.00	40.90	8
6177	OWO WAT V 735	20.992	87.687	45.698	1.00	40.19	8
6178	OWO WAT V 736	46.650	76.726	54.492	1.00	40.47	8
6179	OWO WAT V 737	7.370	47.777	49.173	1.00	41.37	8
6180	OWO WAT V 738	14.279	71.404	55.742	1.00	39.06	8
6181	OWO WAT V 739	13.276	62.067	63.468	1.00	38.95	8
6182	OWO WAT V 740	47.469	55.262	33.788	1.00	41.14	8
6183	OWO WAT V 741	35.182	72.963	73.409	1.00	38.69	8
6184	OWO WAT V 742	28.993	42.749	41.796	1.00	40.22	8
6185	OWO WAT V 743	12.430	51.804	37.400	1.00	39.68	8

No.	Code							
6186	OWO WAT V	744	5.325	85.408	18.950	1.00	41.26	8
6187	OWO WAT V	745	41.721	66.169	73.677	1.00	39.43	8
6188	OWO WAT V	746	16.983	83.519	10.040	1.00	40.83	8
6189	OWO WAT V	747	-3.245	67.552	51.568	1.00	41.14	8
6190	OWO WAT V	748	18.547	72.699	54.390	1.00	40.80	8
6191	OWO WAT V	749	50.952	65.503	10.665	1.00	39.71	8
6192	OWO WAT V	750	6.873	59.618	31.097	1.00	40.61	8
6193	OWO WAT V	751	44.643	72.249	41.540	1.00	40.21	8
6194	OWO WAT V	752	10.333	59.620	21.158	1.00	40.03	8
6195	OWO WAT V	753	25.342	99.843	20.985	1.00	40.88	8
6196	OWO WAT V	754	13.886	67.341	51.913	1.00	41.40	8
6197	OWO WAT V	755	20.669	50.466	56.076	1.00	40.93	8
6198	OWO WAT V	756	16.062	51.352	42.439	0.00	41.08	8
6199	OWO WAT V	757	-2.323	91.095	34.519	1.00	40.69	8
6200	OWO WAT V	758	3.987	63.229	29.353	1.00	41.53	8
6201	OWO WAT V	759	20.694	99.900	18.786	1.00	39.84	8
6202	OWO WAT V	760	21.098	72.115	58.710	1.00	40.04	8
6203	OWO WAT V	761	39.451	41.349	13.513	1.00	40.57	8
6204	OWO WAT V	762	3.185	80.792	42.000	1.00	41.27	8
6205	OWO WAT V	763	15.866	68.494	52.407	1.00	40.90	8
6206	OWO WAT V	764	42.027	79.686	47.257	1.00	40.14	8
6207	OWO WAT V	765	41.063	55.022	68.081	1.00	42.98	8
6208	OWO WAT V	766	15.728	85.261	16.175	1.00	42.08	8
6209	OWO WAT V	767	-8.665	63.541	41.234	1.00	40.98	8
6210	OWO WAT V	768	28.828	85.580	44.440	1.00	40.75	8
6211	OWO WAT V	769	46.504	46.346	43.773	1.00	40.98	8
6212	OWO WAT V	770	52.964	61.429	54.373	1.00	41.49	8
6213	OWO WAT V	771	25.001	95.005	39.507	1.00	41.35	8
6214	OWO WAT V	772	4.818	62.633	59.449	1.00	41.54	8
6215	OWO WAT V	773	18.916	60.160	66.080	1.00	41.09	8
6216	OWO WAT V	774	27.033	88.841	12.019	1.00	41.39	8
6217	OWO WAT V	775	13.803	67.385	57.044	1.00	41.01	8
6218	OWO WAT V	776	11.356	57.839	18.715	1.00	42.40	8
6219	OWO WAT V	777	50.379	81.288	57.286	1.00	38.80	8
6220	OWO WAT V	778	9.739	45.470	44.991	1.00	38.14	8
6221	OWO WAT V	779	26.770	61.900	44.155	1.00	39.31	8
6222	OWO WAT V	780	51.219	53.194	12.424	1.00	42.17	8
6223	OWO WAT V	781	44.665	57.891	62.632	1.00	40.09	8
6224	OWO WAT V	782	-4.332	63.995	45.992	1.00	43.38	8
6225	OWO WAT V	783	54.313	56.319	27.513	1.00	40.34	8
6226	OWO WAT V	784	26.835		8.446	1.00	41.42	8
6227	OWO WAT V	785	53.198	59.584	25.852	1.00	42.12	8

No.	Code							
6228	OWO WAT V	786	19.476	44.845	23.558	1.00	42.02	8
6229	OWO WAT V	787	21.294	88.109	18.755	1.00	42.02	8
6230	OWO WAT V	788	46.461	57.071	56.124	1.00	41.88	8
6231	OWO WAT V	789	52.975	62.124	50.393	1.00	41.59	8
6232	OWO WAT V	790	8.480	108.235	24.221	1.00	42.16	8
6233	OWO WAT V	791	18.517	78.200	51.474	1.00	42.31	8
6234	OWO WAT V	792	-5.150	51.726	49.471	1.00	40.85	8
6235	OWO WAT V	793	36.849	55.447	72.633	1.00	42.13	8
6236	OWO WAT V	794	37.563	79.694	42.494	1.00	41.67	8
6237	OWO WAT V	795	18.818	84.285	16.222	1.00	43.03	8
6238	OWO WAT V	796	36.451	82.759	13.933	1.00	41.71	8
6239	OWO WAT V	797	22.906	51.457	67.808	1.00	41.33	8
6240	OWO WAT V	798	1.528	102.171	36.053	1.00	41.41	8
6241	OWO WAT V	799	36.684	45.294	51.188	1.00	42.88	8
6242	OWO WAT V	800	14.017	67.414	9.706	1.00	41.87	8
6243	OWO WAT V	801	47.832	66.788	18.650	1.00	43.94	8
6244	OWO WAT V	802	34.436	88.324	36.102	1.00	43.03	8
6245	OWO WAT V	803	24.260	67.815	69.741	1.00	42.57	8
6246	OWO WAT V	804	13.129	74.685	52.293	1.00	42.14	8
6247	OWO WAT V	805	16.572	61.433	12.722	1.00	41.61	8
6248	OWO WAT V	806	46.827	51.429	47.524	1.00	42.73	8
6249	OWO WAT V	807	26.546	48.730	71.404	1.00	42.47	8
6250	OWO WAT V	808	33.308	76.572	61.813	1.00	42.82	8
6251	OWO WAT V	809	16.314	98.564	17.363	1.00	42.43	8
6252	OWO WAT V	810	47.038	73.620	61.642	1.00	43.62	8
6253	OWO WAT V	811	23.301	74.541	1.386	1.00	44.52	8
6254	OWO WAT V	812	23.719	88.481	37.683	1.00	40.46	8
6255	OWO WAT V	813	31.440	60.185	76.751	1.00	43.11	8
6256	OWO WAT V	814	16.573	101.550	19.316	1.00	43.14	8
6257	OWO WAT V	815	4.201	55.069	57.571	1.00	42.80	8
6258	OWO WAT V	816	-5.755	72.609	47.268	1.00	41.92	8
6259	OWO WAT V	817	23.802	106.033	28.040	1.00	42.13	8
6260	OWO WAT V	818	28.580	91.427	35.569	1.00	44.77	8
6261	OWO WAT V	819	-6.927	74.495	44.783	1.00	42.39	8
6262	OWO WAT V	820	42.898	80.745	21.282	1.00	42.72	8
6263	OWO WAT V	821	44.594	48.996	19.414	1.00	42.65	8
6264	OWO WAT V	822	22.357	66.608	68.417	1.00	43.63	8
6265	OWO WAT V	823	25.576	50.604	10.991	1.00	43.85	8
6266	OWO WAT V	824	47.220	77.741	25.726	1.00	44.91	8
6267	OWO WAT V	825	16.600	55.517	47.156	1.00	44.33	8
6268	OWO WAT V	826	1.892	50.074	39.116	1.00	42.83	8
6269	OWO WAT V	827	24.440	89.039	35.388	1.00	43.51	8

Index	Label						
6270	OWO WAT V 828	21.480	53.418	43.918	1.00	43.38	8
6271	OWO WAT V 829	14.604	82.502	11.170	1.00	41.84	8
6272	OWO WAT V 830	14.875	62.765	14.779	1.00	42.12	8
6273	OWO WAT V 831	3.701	80.451	16.127	1.00	43.54	8
6274	OWO WAT V 832	48.724	69.957	34.245	1.00	43.72	8
6275	OWO WAT V 833	33.297	59.385	78.143	1.00	45.08	8
6276	OWO WAT V 834	17.544	53.593	51.513	1.00	46.32	8
6277	OWO WAT V 835	12.009	85.638	47.505	1.00	44.72	8
6278	OWO WAT V 836	12.936	54.854	19.146	1.00	42.89	8
6279	OWO WAT V 837	24.005	43.446	16.058	1.00	42.88	8
6280	OWO WAT V 838	2.563	57.015	61.406	1.00	42.80	8
6281	OWO WAT V 839	17.292	81.361	47.327	1.00	43.05	8
6282	OWO WAT V 840	26.973	97.353	24.292	1.00	42.44	8
6283	OWO WAT V 841	24.520	83.741	45.252	1.00	44.48	8
6284	OWO WAT V 842	40.943	89.452	52.148	1.00	43.11	8
6285	OWO WAT V 843	17.748	68.046	54.218	1.00	43.50	8
6286	OWO WAT V 844	53.072	63.513	23.413	1.00	44.03	8
6287	OWO WAT V 845	-1.018	97.468	34.805	1.00	44.47	8
6288	OWO WAT V 846	14.885	69.989	10.290	1.00	44.92	8
6289	OWO WAT V 847	15.800	50.802	20.677	1.00	46.01	8
6290	OWO WAT V 848	25.982	102.118	29.899	1.00	44.94	8
6291	OWO WAT V 849	0.247	52.205	55.531	1.00	45.62	8
6292	OWO WAT V 850	18.884	61.154	6.132	1.00	44.78	8
6293	OWO WAT V 851	11.959	71.909	12.344	1.00	44.89	8
6294	OWO WAT V 852	-1.246	74.512	28.721	1.00	43.38	8
6295	OWO WAT V 853	39.883	48.418	12.157	1.00	44.97	8
6296	OWO WAT V 854	25.484	65.082	75.271	1.00	40.98	8
6297	OWO WAT V 855	31.943	70.931	74.851	1.00	43.51	8
6298	OWO WAT V 856	35.277	49.445	9.697	1.00	45.37	8
6299	OWO WAT V 857	51.902	64.035	51.472	1.00	44.46	8
6300	OWO WAT V 858	5.381	94.650	19.292	1.00	45.20	8
6301	OWO WAT V 859	14.161	52.807	19.941	1.00	44.99	8
6302	OWO WAT V 860	0.976	49.582	46.683	1.00	45.54	8
6303	OWO WAT V 861	11.378	56.649	26.955	1.00	46.98	8
6304	OWO WAT V 862	23.725	48.499	52.657	1.00	43.58	8
6305	OWO WAT V 863	31.619	53.767	4.982	1.00	44.03	8
6306	OWO WAT V 864	21.564	85.158	16.121	1.00	46.34	8
6307	OWO WAT V 865	48.454	57.833	31.207	1.00	45.05	8
6308	OWO WAT V 866	13.820	78.070	11.134	1.00	47.52	8
6309	OWO WAT V 867	53.268	61.611	27.715	1.00	45.47	8
6310	OWO WAT V 868	40.661	42.455	21.732	1.00	43.57	8
6311	OWO WAT V 869	24.881	97.609	17.754	1.00	41.56	8
6312	OWO WAT V 870	46.779	45.863	33.183	1.00	38.06	8
6313	OWO WAT V 871	28.449	47.238	49.410	1.00	44.98	8
6314	OWO WAT V 872	31.691	90.618	16.034	1.00	45.59	8
6315	OWO WAT V 873	45.269	73.271	25.041	1.00	46.87	8
6316	OWO WAT V 874	18.096	50.245	53.261	1.00	45.09	8
6317	OWO WAT V 875	-5.062	64.679	39.520	1.00	44.00	8
6318	OWO WAT V 876	17.006	54.092	49.423	1.00	46.50	8
6319	OWO WAT V 877	39.580	64.107	77.580	1.00	45.16	8
6320	OWO WAT V 878	16.405	56.330	66.595	1.00	44.67	8
6321	OWO WAT V 879	28.253	74.829	0.308	1.00	44.14	8
6322	OWO WAT V 880	26.165	96.027	33.730	1.00	45.73	8
6323	OWO WAT V 881	25.024	94.673	31.684	1.00	45.30	8
6324	OWO WAT V 882	6.382	101.590	23.078	1.00	45.88	8
6325	OWO WAT V 883	20.784	106.032	25.469	1.00	44.92	8
6326	OWO WAT V 884	15.678	70.001	59.741	1.00	47.72	8
6327	OWO WAT V 885	43.426	78.175	61.043	1.00	46.23	8
6328	OWO WAT V 886	10.651	50.358	43.666	1.00	45.79	8
6329	OWO WAT V 887	0.918	49.811	54.685	1.00	45.38	8
6330	OWO WAT V 888	24.371	106.460	30.607	1.00	46.20	8
6331	OWO WAT V 889	7.425	93.590	16.782	1.00	45.79	8
6332	OWO WAT V 890	20.314	66.093	66.469	1.00	48.35	8
6333	OWO WAT V 891	12.244	93.829	44.935	1.00	47.03	8
6334	OWO WAT V 892	2.847	56.085	32.137	1.00	48.09	8
6335	OWO WAT V 893	15.592	72.473	53.588	1.00	46.24	8
6336	OWO WAT V 894	45.552	49.673	29.081	1.00	46.73	8
6337	OWO WAT V 895	48.867	72.566	56.249	1.00	45.76	8
6338	OWO WAT V 896	51.123	56.932	37.786	1.00	40.59	8
6339	OWO WAT V 897	30.394	49.733	9.142	1.00	48.10	8
6340	OWO WAT V 898	27.101	98.920	22.257	1.00	46.90	8
6341	OWO WAT V 899	14.712	51.617	24.308	1.00	47.17	8
6342	OWO WAT V 900	46.781	50.533	27.214	1.00	47.00	8
6343	OWO WAT W 1	-6.114	65.250	32.669	1.00	47.96	8
6344	OWO WAT W 2	47.645	71.480	60.598	1.00	45.49	8
6345	OWO WAT W 3	48.802	54.151	54.398	1.00	46.37	8
6346	OWO WAT W 4	4.605	105.268	26.990	1.00	46.92	8
6347	OWO WAT W 5	42.143	73.136	41.061	1.00	45.02	8
6348	OWO WAT W 6	26.105	89.154	41.563	1.00	47.90	8
6349	OWO WAT W 7	3.202	48.853	55.140	1.00	47.23	8
6350	OWO WAT W 8	-2.318	61.724	33.458	1.00	48.00	8
6351	OWO WAT W 9	38.569	50.803	7.204	1.00	46.95	8
6352	OWO WAT W 10	2.160	103.336	25.390	1.00	47.08	8
6353	OWO WAT W 11	27.464	41.903	36.189	1.00	42.96	8

6396	OWO	WAT	W	54	49.464	48.219	14.354	1.00	55.58	8
6397	OWO	WAT	W	55	-2.823	49.394	48.028	1.00	50.11	8
6398	OWO	WAT	W	56	22.263	105.611	23.242	1.00	50.89	8
6399	OWO	WAT	W	57	16.395	51.479	56.998	1.00	54.48	8
6400	OWO	WAT	W	58	46.461	51.028	12.082	1.00	51.52	8
6401	OWO	WAT	W	59	35.996	61.651	5.602	1.00	53.96	8
6402	OWO	WAT	W	60	25.277	46.888	50.958	1.00	53.30	8
6403	OWO	WAT	W	61	15.763	94.604	17.487	1.00	49.77	8
6404	OWO	WAT	W	62	42.937	61.534	7.934	1.00	47.74	8
6405	OWO	WAT	W	63	24.840	46.516	44.072	1.00	51.69	8
6406	OWO	WAT	W	64	29.111	92.028	40.197	1.00	48.35	8
6407	OWO	WAT	W	65	4.689	89.221	19.393	1.00	50.44	8
6408	OWO	WAT	W	66	11.456	93.383	15.046	1.00	58.07	8
6409	OWO	WAT	W	67	15.227	108.500	35.816	1.00	52.43	8
6410	OWO	WAT	W	68	42.860	74.904	32.815	1.00	52.70	8
6411	OWO	WAT	W	69	48.829	69.006	23.605	1.00	55.19	8
6412	OWO	WAT	W	70	15.485	68.666	11.962	1.00	52.48	8
6413	OWO	WAT	W	71	-5.047	98.836	30.411	1.00	52.33	8
6414	OWO	WAT	W	72	38.899	69.928	3.024	1.00	53.43	8
6415	OWO	WAT	W	73	47.563	47.253	16.948	1.00	53.16	8
6416	OWO	WAT	W	74	8.621	51.721	57.810	1.00	51.63	8
6417	OWO	WAT	W	75	52.887	59.020	49.147	1.00	50.28	8
6418	OWO	WAT	W	76	4.037	61.529	30.846	1.00	54.62	8
6419	OWO	WAT	W	77	4.332	83.232	16.091	1.00	55.82	8
6420	OWO	WAT	W	78	36.037	53.977	74.978	1.00	55.15	8
6421	OWO	WAT	W	79	51.512	53.978	47.701	0.00	54.18	8
6422	OWO	WAT	W	80	14.797	91.920	17.273	1.00	51.87	8
6423	OWO	WAT	W	81	19.801	84.369	8.356	1.00	55.63	8
6424	OWO	WAT	W	82	31.496	46.252	52.107	1.00	55.15	8
6425	OWO	WAT	W	83	-1.823	79.829	19.762	1.00	53.76	8
6426	OWO	WAT	W	84	29.256	92.834	20.731	1.00	53.49	8
6427	OWO	WAT	W	85	-1.824	76.279	32.563	1.00	55.76	8
6428	OWO	WAT	W	86	32.061	46.088	49.588	1.00	55.95	8
6429	OWO	WAT	W	87	41.327	42.439	33.669	1.00	55.86	8
6430	OWO	WAT	W	88	43.153	80.324	3.532	1.00	52.67	8
6431	OWO	WAT	W	89	47.799	76.801	61.841	1.00	55.93	8
6432	OWO	WAT	W	90	35.708	85.333	37.348	1.00	54.64	8
6433	OWO	WAT	W	91	52.110	65.452	56.038	1.00	53.77	8
6434	OWO	WAT	W	92	16.996	79.167	47.458	1.00	56.65	8
6435	OWO	WAT	W	93	27.626	49.502	11.515	1.00	60.46	8
6436	OWO	WAT	W	94	29.317	47.547	14.906	1.00	61.23	8
6437	OWO	WAT	W	95	51.306	65.455	30.794	0.00	67.31	8

6354	OWO	WAT	W	12	49.209	50.152	22.583	1.00	48.25	8
6355	OWO	WAT	W	13	10.266	57.487	62.131	1.00	43.39	8
6356	OWO	WAT	W	14	54.127	66.185	63.244	1.00	49.78	8
6357	OWO	WAT	W	15	19.083	49.330	63.179	1.00	49.30	8
6358	OWO	WAT	W	16	39.385	74.287	38.793	1.00	46.50	8
6359	OWO	WAT	W	17	-1.255	52.062	41.948	1.00	49.09	8
6360	OWO	WAT	W	18	16.630	69.708	8.569	1.00	48.63	8
6361	OWO	WAT	W	19	29.289	54.314	9.208	1.00	49.09	8
6362	OWO	WAT	W	20	38.590	83.950	31.916	1.00	48.14	8
6363	OWO	WAT	W	21	47.212	44.923	37.558	1.00	49.07	8
6364	OWO	WAT	W	22	2.016	104.557	33.998	1.00	50.31	8
6365	OWO	WAT	W	23	5.910	79.410	15.810	1.00	46.74	8
6366	OWO	WAT	W	24	18.824	46.920	16.693	1.00	44.54	8
6367	OWO	WAT	W	25	21.253	90.457	18.335	1.00	47.12	8
6368	OWO	WAT	W	26	46.674	72.576	34.746	1.00	48.69	8
6369	OWO	WAT	W	27	17.804	54.470	14.884	1.00	50.05	8
6370	OWO	WAT	W	28	-3.319	64.779	54.655	1.00	49.95	8
6371	OWO	WAT	W	29	20.035	59.810	69.328	1.00	47.36	8
6372	OWO	WAT	W	30	6.598	98.214	19.529	1.00	50.29	8
6373	OWO	WAT	W	31	29.388	57.634	5.275	1.00	49.23	8
6374	OWO	WAT	W	32	25.881	83.371	49.453	1.00	49.89	8
6375	OWO	WAT	W	33	46.828	48.534	13.616	1.00	48.17	8
6376	OWO	WAT	W	34	52.912	62.999	14.852	1.00	46.56	8
6377	OWO	WAT	W	35	2.107	83.103	18.070	1.00	48.10	8
6378	OWO	WAT	W	36	41.486	67.797	7.736	1.00	50.23	8
6379	OWO	WAT	W	37	9.492	94.962	15.817	1.00	48.00	8
6380	OWO	WAT	W	38	28.348	95.134	21.001	1.00	48.19	8
6381	OWO	WAT	W	39	32.401	92.434	25.439	1.00	50.58	8
6382	OWO	WAT	W	40	46.035	46.956	48.050	1.00	50.92	8
6383	OWO	WAT	W	41	34.909	51.572	71.159	1.00	49.38	8
6384	OWO	WAT	W	42	29.114	76.486	61.672	1.00	49.83	8
6385	OWO	WAT	W	43	-3.044	91.942	40.975	1.00	51.92	8
6386	OWO	WAT	W	44	-3.395	74.451	28.328	1.00	46.84	8
6387	OWO	WAT	W	45	20.187	89.421	44.602	1.00	49.88	8
6388	OWO	WAT	W	46	-4.080	54.699	41.545	1.00	50.13	8
6389	OWO	WAT	W	47	38.425	87.733	31.621	1.00	49.47	8
6390	OWO	WAT	W	48	4.305	62.413	62.323	1.00	49.60	8
6391	OWO	WAT	W	49	6.229	89.367	44.095	1.00	53.06	8
6392	OWO	WAT	W	50	49.747	66.902	23.155	0.00	51.47	8
6393	OWO	WAT	W	51	40.855	78.417	26.292	1.00	48.79	8
6394	OWO	WAT	W	52	17.489	50.714	66.895	1.00	52.19	8
6395	OWO	WAT	W	53	34.327	92.622	27.782	1.00	49.40	8

ID	Label	Index						
6438	OWO WAT W	96	21.436	52.799	12.475	1.00	60.39	8
6439	OWO WAT W	97	24.857	50.122	68.228	1.00	61.97	8
6440	OWO WAT W	98	53.436	60.942	47.809	1.00	59.88	8
6441	OWO WAT W	99	26.545	99.863	28.613	1.00	66.13	8
6442	OWO WAT W	100	28.187	94.100	34.809	1.00	47.55	8
6443	OWO WAT W	101	46.501	68.477	9.327	1.00	62.49	8
6444	OWO WAT W	102	41.335	80.622	32.546	1.00	56.46	8
6445	OWO WAT W	103	49.090	47.019	29.937	1.00	51.60	8
6446	OWO WAT W	104	29.677	70.505	75.480	1.00	55.52	8
6447	OWO WAT W	105	10.580	70.552	56.020	1.00	45.60	8
6448	OWO WAT W	106	-5.437	61.460	36.195	1.00	49.58	8
6449	OWO WAT W	107	41.636	42.378	28.372	1.00	47.73	8
6450	OWO WAT W	108	48.134	51.375	29.584	1.00	53.67	8
6451	OWO WAT W	109	20.029	46.534	41.141	1.00	45.69	8
6452	OWO WAT W	110	39.076	61.857	77.827	1.00	61.24	8
6453	OWO WAT W	111	40.140	79.602	28.296	1.00	58.08	8
6454	OWO WAT W	112	24.479	41.686	38.003	1.00	48.71	8
6455	OWO WAT W	113	18.748	86.522	18.736	1.00	54.95	8
6456	OWO WAT W	114	26.670	86.155	43.878	1.00	45.83	8
6457	OWO WAT W	115	34.014	44.101	58.947	1.00	55.80	8
6458	OWO WAT W	116	44.085	44.358	32.579	1.00	66.42	8
6459	OWO WAT W	117	2.549	102.526	41.260	1.00	56.37	8
6460	OWO WAT W	118	10.042	59.115	64.251	1.00	53.53	8
6461	OWO WAT W	119	52.498	59.271	37.676	1.00	46.44	8
6462	OWO WAT W	120	49.412	68.479	30.738	0.00	48.32	8
6463	OWO WAT W	121	39.604	81.174	29.899	1.00	40.72	8
6464	OWO WAT W	122	52.578	62.726	30.463	0.00	55.27	8
6465	OWO WAT W	123	32.284	38.992	32.423	1.00	44.19	8
6466	OWO WAT W	124	54.342	58.298	13.900	1.00	48.84	8
6467	OWO WAT W	125	53.831	60.018	17.609	1.00	53.12	8
6468	OWO WAT W	126	37.548	48.910	67.790	1.00	63.29	8
6469	OWO WAT W	127	16.364	67.201	64.210	1.00	54.77	8
6470	OWO WAT W	128	35.507	88.726	18.930	1.00	44.18	8
6471	OWO WAT W	129	49.585	56.011	59.240	1.00	44.54	8
6472	OWO WAT W	130	13.470	54.095	28.765	1.00	39.17	8
6473	OWO WAT W	131	11.141	90.680	16.198	1.00	49.46	8
6474	OWO WAT W	132	-9.184	65.150	46.616	1.00	52.99	8
6475	OWO WAT W	133	44.910	67.182	7.757	1.00	52.67	8
6476	OWO WAT W	134	38.968	68.370	71.252	1.00	37.18	8
6477	OWO WAT W	135	9.962	106.661	35.598	1.00	48.81	8
6478	OWO WAT W	136	-0.018	83.329	37.562	1.00	49.93	8
6479	OWO WAT W	137	6.077	95.625	42.890	1.00	41.31	8
6480	OWO WAT W	138	38.033	48.935	8.928	1.00	56.21	8
6481	OWO WAT W	139	22.720	48.167	43.406	1.00	43.05	8
6482	OWO WAT W	140	16.160	51.440	42.398	1.00	42.33	8
6483	OWO WAT W	141	51.286	65.520	30.747	0.00	58.93	8
6484	OWO WAT W	142	22.870	83.783	-0.279	1.00	39.91	8
6485	OWO WAT W	143	23.492	85.411	14.742	1.00	45.55	8
6486	OWO WAT W	144	30.609	38.135	34.869	1.00	53.23	8
6487	OWO WAT W	145	51.546	53.971	47.725	0.00	54.90	8
6488	OWO WAT W	146	37.344	40.493	33.234	1.00	46.02	8
6489	OWO WAT W	147	35.805	47.572	62.190	1.00	53.24	8
6490	OWO WAT W	148	32.439	62.293	76.111	1.00	52.51	8
6491	OWO WAT W	149	24.077	90.751	37.673	1.00	48.00	8
6492	OWO WAT W	150	20.655	50.869	68.464	1.00	51.34	8
6493	OWO WAT W	151	42.359	76.800	31.368	1.00	54.68	8
6494	OWO WAT W	152	40.991	84.869	28.522	1.00	51.94	8
6495	OWO WAT W	153	-3.448	59.486	34.298	1.00	60.14	8
6496	OWO WAT W	154	24.275	50.044	65.629	1.00	48.99	8
6497	OWO WAT W	155	24.898	47.635	57.042	1.00	50.00	8
6498	OWO WAT W	156	46.911	73.376	32.086	1.00	65.97	8
6499	OWO WAT W	157	12.448	63.643	14.806	1.00	50.68	8
6500	OWO WAT W	158	17.367	83.516	7.750	1.00	59.03	8
6501	OWO WAT W	159	38.537	87.429	19.244	1.00	48.05	8
6502	OWO WAT W	160	49.397	68.482	30.753	0.00	47.12	8
6503	OWO WAT W	161	52.562	62.737	30.460	0.00	55.04	8
6504	OWO WAT W	162	17.100	92.730	17.506	1.00	48.89	8
6505	OWO WAT W	163	54.143	65.666	53.272	1.00	49.78	8
6506	OWO WAT W	164	35.140	89.217	21.335	0.00	49.55	8
6507	OWO WAT W	165	40.864	85.144	25.201	1.00	51.96	8
6508	OWO WAT W	166	0.129	71.062	53.859	1.00	50.46	8
6509	OWO WAT W	167	19.749	95.732	16.395	1.00	54.02	8
6510	OWO WAT W	168	45.089	55.696	66.763	1.00	48.69	8
6511	OWO WAT W	169	29.920	93.952	27.969	1.00	57.82	8
6512	OWO WAT W	170	-1.140	103.281	29.038	1.00	59.68	8
6513	OWO WAT W	171	0.493	67.332	27.261	1.00	46.75	8
6514	OWO WAT W	172	11.663	49.273	52.961	1.00	47.65	8
6515	OWO WAT W	173	19.395	43.670	27.526	0.00	52.64	8
6516	OWO WAT W	174	12.558	73.429	10.221	1.00	52.09	8
6517	OWO WAT W	175	47.725	72.168	25.475	1.00	55.03	8
6518	OWO WAT W	176	37.354	46.525	53.992	1.00	62.86	8
6519	OWO WAT W	177	6.566	77.689	48.060	1.00	54.04	8
6520	OWO WAT W	178	27.239	80.756	54.362	1.00	54.49	8
6521	OWO WAT W	179	29.136	79.115	59.854	1.00	51.47	8

6522	OWO	WAT	W	180	51.301	65.512	30.778	1.00	58.87	8
6523	OWO	WAT	W	181	49.729	67.077	23.232	1.00	50.64	8
6524	OWO	WAT	W	182	15.428	63.127	65.725	1.00	53.21	8
6525	OWO	WAT	W	183	28.316	47.473	59.001	1.00	52.58	8
6526	OWO	WAT	W	184	11.167	50.777	21.789	1.00	55.21	8
6527	OWO	WAT	W	185	39.667	45.779	11.484	1.00	50.46	8
6528	OWO	WAT	W	186	9.302	78.481	48.408	1.00	50.14	8
6529	OWO	WAT	W	187	-2.511	95.777	25.082	1.00	50.26	8
6530	OWO	WAT	W	188	49.525	52.221	41.616	1.00	53.78	8
6531	OWO	WAT	W	189	33.219	90.787	32.195	1.00	53.22	8
6532	OWO	WAT	W	190	18.629	63.971	67.355	1.00	54.97	8
6533	OWO	WAT	W	191	6.996	55.402	34.790	1.00	49.49	8
6534	OWO	WAT	W	192	50.269	70.807	60.070	1.00	54.80	8
6535	OWO	WAT	W	193	-3.948	89.479	31.518	0.00	54.80	8
6536	OWO	WAT	W	194	4.036	51.795	57.529	1.00	56.49	8
6537	OWO	WAT	W	195	15.790	48.084	23.036	1.00	52.39	8
6538	OWO	WAT	W	196	22.577	105.708	45.361	1.00	51.14	8
6539	OWO	WAT	W	197	7.453	103.921	23.508	1.00	50.15	8
6540	OWO	WAT	W	198	37.899	80.588	45.384	1.00	52.20	8
6541	OWO	WAT	W	199	19.774	74.879	53.890	1.00	55.07	8
6542	OWO	WAT	W	200	50.055	68.873	30.956	1.00	30.00	8
6543	OWO	WAT	W	201	53.330	63.201	30.956	1.00	37.00	8
6544	OWO	WAT	W	202	35.086	89.130	21.284	1.00	49.00	8
6545	OWO	WAT	W	203	2.339	51.858	35.796	1.00	50.00	8
6546	OWO	WAT	W	204	19.180	43.755	27.572	1.00	50.00	8
6547	OWO	WAT	W	205	51.693	55.504	46.921	1.00	51.00	8
6548	OWO	WAT	W	206	31.811	80.217	54.661	1.00	51.00	8
6549	OWO	WAT	W	207	11.695	77.786	12.093	1.00	51.00	8
6550	OWO	WAT	W	208	29.940	46.996	53.693	1.00	52.00	8
6551	OWO	WAT	W	209	7.251	102.500	40.633	1.00	52.00	8
6552	OWO	WAT	W	210	23.858	91.561	17.414	1.00	52.00	8
6553	OWO	WAT	W	211	6.783	49.832	40.633	1.00	52.00	8
6554	OWO	WAT	W	212	44.910	47.806	22.735	1.00	52.00	8
6555	OWO	WAT	W	213	36.255	46.591	10.158	1.00	52.00	8
6556	OWO	WAT	W	214	27.601	61.581	77.880	1.00	52.00	8
6557	OWO	WAT	W	215	27.133	98.043	33.861	1.00	53.00	8
6558	OWO	WAT	W	216	18.479	55.504	45.470	1.00	53.00	8
6559	OWO	WAT	W	217	9.122	47.401	46.438	1.00	53.00	8
6560	OWO	WAT	W	218	9.590	66.037	16.447	1.00	53.00	8
6561	OWO	WAT	W	219	13.333	91.966	46.438	1.00	53.00	8

122

CLAIMS:

1.   A method of producing a variant of a parent maltogenic alpha-amylase, said method comprising

      a)   modeling the parent alpha-amylase on the three-dimensional structure of SEQ ID NO: 1 depicted in the Appendix to produce a three-dimensional structure of the parent alpha-amylase;

      b)   identifying in the three-dimensional structure obtained in step (a) at least one structural part of the parent wherein an alteration in said structural part is predicted to result in said altered property;

      c)   modifying the sequence of a nucleic acid encoding the parent alpha-amylase to produce a nucleic acid encoding a deletion, insertion, or substitution of one or more amino acids at a position corresponding to said structural part; and

      d)   expressing the modified nucleic acid in a host cell to produce the variant alpha-amylase,

wherein the variant has alpha-amylase enzymatic activity and has at least one altered property relative to the parent.

2.   The method of claim 1, wherein the altered property is pH dependent activity, thermostability, substrate cleavage pattern, specific activity of cleavage, transglycosylation, ability to reduce retrogradation of starch, ability to reduce staling of bread, substrate specificity, substrate binding or calcium binding.

3.   A method of constructing a variant of a parent maltogenic alpha-amylase, which method comprises:

      a)   identifying an amino acid residue which is within 15 Å (in particular 10 Å) from an active site residue of the parent amylase in the three-dimensional structure of said parent, and which is involved in electrostatic or hydrophobic interactions with an active site residue;

      b)   substituting said amino acid residue with another amino acid residue which changes the electrostatic and/or hydrophobic surroundings of an active site residue, and which can be accommodated in the structure;

      c)   optionally repeating steps a) and b) recursively;

      d)   optionally, making alterations each of which is an insertion, a deletion or a substitution of an amino acid residue at one or mor   positions other than b),

123

e)   preparing the variant resulting from steps a) - d);

f)   testing the pH dependent activity of said variant; and

g)   optionally repeating steps a) - f) recursively; and

h)   selecting a variant having an altered pH dependent activity as compared

5   to the parent amylase.

4.   A method of constructing a variant of a parent maltogenic alpha-amylase, which method comprises:

a)   identifying an internal cavity or crevice in the three-dimensional structure of said parent;

10   b)   substituting an amino acid residue in the neighborhood of the cavity or crevice with another amino acid residue which increases the hydrophobic interaction and/or fills out or reduces the size of the cavity or crevice;

c)   optionally repeating steps a) and b) recursively;

d)   optionally, making alterations each of which is an insertion, a deletion or a

15   substitution of an amino acid residue at one or more positions other than b),

e)   preparing the variant resulting from steps a) - d);

f)   testing the thermostability of said variant; and

g)   optionally repeating steps a) - f) recursively; and

h)   selecting a variant having increased thermostability as compared to the

20   parent amylase.

5.   The method of claim 4, wherein the substitution of the amino acid residue results in increasing the hydrophobic interaction, substitution with proline, substitution of histidine with another amino acid, stabilization of calcium binding, introduction of an interdomain disulfide bond, removal of a deamidation site, altering a hydrogen

25   bond contact, filling in an internal structural cavity with an amino acid with a bulkier side group, introduction of interdomain interactions, altering charge distribution, helix capping, or introduction of a salt bridge.

6.   A method of constructing a variant of a parent maltogenic alpha-amylase, which method comprises:

30   a)   identifying an amino acid residue within 10 Å from a calcium binding site in the three dimensional structure of the amylase;

b)   substituting the amino acid residue with another amino acid residue so as to improve the interaction with the calcium ion;

c)   optionally rep ating steps a) and b) recursively;

124

d)   optionally, making alterations each of which is an insertion, a deletion or a substitution of an amino acid residue at one or more positions other than b),

e)   preparing the variant resulting from steps a) - d);

f)   testing the thermostability of said variant; and

g)   optionally repeating steps a) - f) recursively; and

h)   selecting a variant having increased thermostability as compared to the parent amylase.

7.   A method of constructing a variant of a parent maltogenic alpha-amylase, which method comprises:

a)   identifying the substrate binding area in a model of the three-dimensional structure of the parent amylase;

b)   modifying the substrate binding area by an amino acid substitution, deletion or insertion;

c)   optionally repeating step b) recursively;

d)   optionally, making alterations each of which is an insertion, a deletion or a substitution of an amino acid residue at one or more positions other than b),

e)   preparing the variant resulting from steps a) - d);

f)   testing the substrate-cleavage pattern of the variant.

g)   optionally repeating steps a) - f) recursively; and

h)   selecting a variant having an altered substrate-cleavage pattern as compared to the parent amylase.

8.   A method for producing a maltogenic alpha-amylase variant, comprising:

a)   constructing the variant by the method of any of claims 2-7;

b)   transforming a microorganism with a DNA sequence encoding the variant;

c)   cultivating the transformed microorganism under conditions which are conducive for producing the variant, and

d)   optionally, recovering the variant from the resulting culture broth.

9.   A polypeptide which:

a)   has maltogenic alpha-amylase activity;

b)   has at least 70 % identity to SEQ ID NO: 1;

c)   comprises an amino acid modification compared to SEQ ID NO: 1 at a position corresponding to D127, V129, F188, A229, Y258, V281, F284, T288, N327, M330, G370, N371, and/or D372; and

d)   has altered pH dependent activity as compared to the polypeptide of SEQ ID NO: 1.

10.   The polypeptide of claim 9, wherein the modification comprises a substitution corresponding to D127N/L, V129S/T/G/V, F188E/K/H, A229S/T/G/V, Y258E/D/K/R/F/N, V281L/T, F284K/H/D/E/Y, T288E/K/R, N327D, M330L/F/I/D/E/K, G370N, N371D/E/G/K, and/or D372N/V.

5 11.   A polypeptide which:
       a)   has maltogenic alpha-amylase activity;
       b)   has at least 70 % identity to SEQ ID NO: 1; and
       c)   comprises an amino acid modification compared to SEQ ID NO: 1 at a
       position corresponding to Q13, I16, D17, N26, N28, P29, A30, S32, Y33, G34,
10   L35, K40, M45, P73, V74, D76, N77, D79, N86, R95, N99, I100, H103, Q119,
       N120, N131, S141, T142, A148, N152, A163, H169, N171, G172, I174, N176,
       N187, F188, A192, Q201, N203, H220, N234, G236, Q247, K249, D261, N266,
       L268, R272, N275, N276, V279, N280, V281, D285, N287, F297, Q299, N305,
       K316, N320, L321, N327, A341, N342, A348, Q365, N371, N375, M378, G397,
15   A381, F389, N401, A403, K425, N436, S442, N454, N468, N474, S479, A483,
       A486, V487, S493, T494, S495, A496, S497, A498, Q500, N507, I510, N513,
       K520, Q526, A555, A564, S573, N575, Q581, S583, F586, K589, N595, G618,
       N621, Q624, A629, F636, K645, N664 and/or T681; and
       d)   has improved stability compared to the polypeptide of SEQ ID NO: 1.

20 12.   The polypeptide of claim 11, wherein the modification comprises at a position corresponding to K40, V74, H103, S141, T142, F188, H220, N234, K249, D261, L268, V279, N342, H344, G397, A403, K425, S442, S479, S493, T494, S495, A496, S497, A498, Q500, K520, A555 and/or N595; preferably a substitution corresponding to K40R, V74P, H103Y/V/I/L/F/Y, S141P, T142A, F188I/L, H220Y/L/M, N234P,
25 K249P, D261G, L268P, V279P, N342P, H344E/Q/N/D/Y, G397P, A403P, K425E, S442P, S479P, S493P, T494P, S495P, A496P, S497P, A498P, Q500P, K520R, A555P and/or N595I.

13.   The polypeptide of claim 11 or 12, wherein the modification comprises at a position corresponding to D17, N28, P29, A30, S32, Y33, G34, R95, H103, N131,
30 H169, I174 and/or Q201 such as to improve calcium coordination, preferably a substitution corresponding to D17Q/E, A30D/M/L/A/V/I/E/Q, S32D/E/N/Q, R95M/L/A/V/I/E/Q, H103Y/N/Q/D/E, N131D, H169N/D/E/Q, I174E/Q, Q201E.

14.   The polypeptide of any of claims 11-13, wherein the modification comprises a substitution at a position corresponding to Q13, N26, N77, N86, N99, Q119, N120,

N131, N152, N171, N176, N187, Q201, N203, N234, Q247, N266, N275, N276, N280, N287, Q299, N320, N327, N342, Q365, N371, N375, N401, N436, N454, N468, N474, Q500, N507, N513, Q526, N575, Q581, N621, Q624 and/or N664 such as to remove a deamidation site, preferably a substitution corresponding to

5   Q13S/T/A/V/L/I/F/M, N26S/T/A/V/L/I, N77S/T/A/V/L/I, N86S/T/A/V/L/I, N99T/S/V/L, Q119T/S, N120S/T/A/V/L/I, N131S/T/A/V/L/I, N152T/S/V/L, N171Y/D/S/T, N176S/T/A/V/L/I, N187S/T/A/V/L/I, Q201S/T/A/V/L/I/F/M, N203D/S/T/A/V/L/I, N234S/T/A/V/L/I, Q247S/T/A/V/L/I/F/M, N266S/T/A/V/L/I, N275S/T/A/V/L/I, N276S/T/A/V/L/I, N280S/T/A/V/L/I, N287S/T/A/V/L/I, Q299L/T/S, N320S/T/A/V/L/I,

10 N327S/T/A/V/L/I, N342S/T/A/V/L/I, Q365S/T/A/V/L/I, N371S/T/A/V/L/I, N375S/T/A/V/L/I, N401S/T/A/V/L/I, N436S/T/A/V/L/I, N454D/S/T/A/V/L/I, N468D/S/T/A/V/L/I, N474D/S/T/A/V/L/I, Q500S/T/A/V/L/I/F/M, N507S/T/A/V/L/I, N513S/T/A/V/L/I, Q526 D/S/T/A/V/L/I, N575S/T/A/V/L/I, Q581S/T/A/V/L/I/F/M, N621S/T/A/V/L/I Q624S/T/A/V/L/I/F/M and/or N664D/S/T/A/V/L/I.

15 15.   The polypeptide of any of claims 11-14, wherein the modification comprises a substitution at a position corresponding to I16, L35, M45, P73, D76, D79, A192, I100, A148, A163+G172, L268, V281, D285, L321, F297, N305, K316, S573, A341, M378, A381, F389, A483, A486, I510, A564, F586, K589, F636, K645, A629, and/or T681 such as to improve hydrogen bond contact, preferably a substitution corresponding to

20 I16T/D/N, L35Q, M45K, P73Q, D76E, D79E/Y, A192S/D/N, I100T/S/D/N/E/Q, A148D/N/E/Q/S/T/R/K, A163Y+G172S/D/N, L268R/K, V281/Q, D285R/K, L321Q, F297N/D/Q/E, N305K/R, K316N/D, S573N/D, A341R/K, M378R/K, A381S/D/N, F389Y, A483S/D/N, A486Q/E, I510R/K, A564S/D/N, F586S/D/N, 'K589S/D/Q/N, F636Y, K645T, A629N/D/E/Q, and/or T681D/N/E/Q/S.

25 16.   The polypeptide of any of claims 11-15, wherein the modification comprises substitutions such as to introduce one or more interdomain disulfide bonds, preferably corresponding to G236C + S583C, G618C + R272C, and/or A348C + V487C.

17.   The polypeptide of any of claims 11-16, wherein the substitution at a position corresponding to L51, L75, L78, G88, G91, T94, V114, I125, V126, T134, G157,

30 L217, S235, G236, V254, V279, V281, L286, V289, I290, V308, L321, I325, D326, L343, F349, S353, I359, I405, L448, Q449, L452, I470, G509, V515, S583, G625, L627, L628 and/or A670 so as to fill an internal cavity or crevice, preferably a substitution corresponding to L51W, L75F/Y, L78I, G88A/V/T, G91T/S/V/N, T94V/I/L, V114V/I/L, I125L/M/F/Y/W, V126I/L, T134V/I/L/M/F/Y/W, G157A/V/I/L,

35 L217V/I/M/F/Y/W, S235I/L/M/F/Y/W, G236A/V/I/L/M/F/Y/W, V254I/L/M/F/Y/W,

V279M/I/L/F, V281I/L/M/F/Y/W, L286F, V289I/L/R, I290M/L/F, V308I/L/M/F/Y/W, L321I/M/F/Y/W, I325L/M/F/Y/W, D326E/Q, L343M/F/Y/W, F349W/Y, S353V/I/L, I359L/M/F/Y/W, I405M/L/Y/F/W, L448Y, Q449Y, L452M/Y/F/W, I470M/L/F, G509A/V/I/L/M/S/T/D/N, V515I/L, S583V/I/L/V, G625A/V/I/L/M/F/Y/W, L627M/F/Y,
5 L628M/I/F/Y/W, A670V/I/L/M/F/Y/W and/or L217 in combination with L75 (e.g. L217F/Y in combination with L75F/Y).

18.   The polypeptide of any of claims 11-17, wherein the modification comprises a substitution at a position corresponding to N106, N320 and Q624 so as to create a salt bridge, preferably a substitution corresponding to N106R, N320E/D and/or
10 Q624E.

19.   The polypeptide of any of claims 11-18, wherein the modification comprises a substitution at a position corresponding to K244 and/or K316 such as to alter the charge distribution, preferably a substitution corresponding to K244S and/or K316G/N/D.

15 20.   The polypeptide of any of claims 11-19, wherein the modification comprises a substitution at a position corresponding to V281 and/or A629 such as to alter the binding site, preferably a substitution corresponding to V281Q and/or A629N/D/E/Q.

21.   The polypeptide of any of claims 11-20, wherein the modification comprises substitutions such as to alter the interdomain interaction at a position corresponding
20 to F143+F194+L78, A341+A348+L398+I415+T439+L464+L465, L557, S240+L268, Q208+L628, F427+Q500+N507+M508+S573 and/or I510+V620, preferably substitutions corresponding to F143Y+F194Y+L78Y/F/W/E/Q, A341S/D/N+A348V/I/L+L398E/Q/N/D+I415E/Q+T439D/E/Q/N+L464D/E+L465D/E/N/ Q/R/K, L557Q/E/N/D, S240D/E/N/Q+L268D/E/N/Q/R/K, Q208D/E/Q+L628E/Q/N/D,
25 F427E/Q/R/K/Y+Q500Y+N507Q/E/D+M508K/R/E/Q+S573D/E/N/Q;   and/or I510D/E/N/Q/S+V620D/E/N/Q.

22.   A polypeptide which:
       a)   has maltogenic alpha-amylase activity;
       b)   has at least 70 % identity to SEQ ID NO: 1;
30   c)   comprises an amino acid modification compared to SEQ ID NO: 1 at a position corresponding to P191, A192, G193, F194 and/or S195; and
       d)   has higher specific amylase activity than the polypeptide of SEQ ID NO: 1.

128

23.   The polypeptide of claim 22, wherein the modification comprises a deletion, preferably the deletion Δ (191-195).

24.   The polypeptide of claim 22, wherein the modification comprises insertion, preferably 192-A-193.

5   25.   A polypeptide which:
           a)   has maltogenic alpha-amylase activity;
           b)   has at least 70 % identity to SEQ ID NO: 1;
           c)   comprises an amino acid modification compared to SEQ ID NO: 1 at a position corresponding to A30, K40, N115, T142, F188, T189, P191, A192,
10   G193, F194, S195, D261, T288, N327, K425, K520 and/or N595; and
           d)   has a higher ability than the polypeptide of SEQ ID NO: 1 to reduce retrogradation of starch and/or staling of bread.

26.   The polypeptide of claim 26, wherein the modification comprises A30D, K40R, N115D, T142A F188L, T189Y, Δ (191-195), D261G, T288P, N327S, K425E, K520R
15 and/or N595I.

27.   A nucleic acid sequence encoding the polypeptide of any of claims 9-26, preferably operably linked to one or more control sequences which direct the expression of the variant in a suitable expression host.

28.   A recombinant expression vector comprising the nucleic acid sequence of
20 claim 27, a promoter, and transcriptional and translational stop signals, and preferably further comprising a selectable marker.

29.   A transformed host cell comprising the nucleic acid sequence of claim 27 or the vector of claim 28.

30.   A method for producing the polypeptide of any of claims 9-26, comprising:
25   a)   cultivating the transformed host cell of claim 29 under conditions conducive to expression of the variant; and
           b)   recovering the variant.

31.   A process for preparing a dough or a baked product prepared from the dough which comprises adding the polypeptide of any of claims 9-26, or a variant produced

by the method of any of claims 1-8 to the dough in an amount which is effective to retard the staling of the bread.

32.   The process of claim 31, wherein the variant is added in an amount of 0.1-5 mg per kg of flour, preferably 0.5-2 mg/kg.
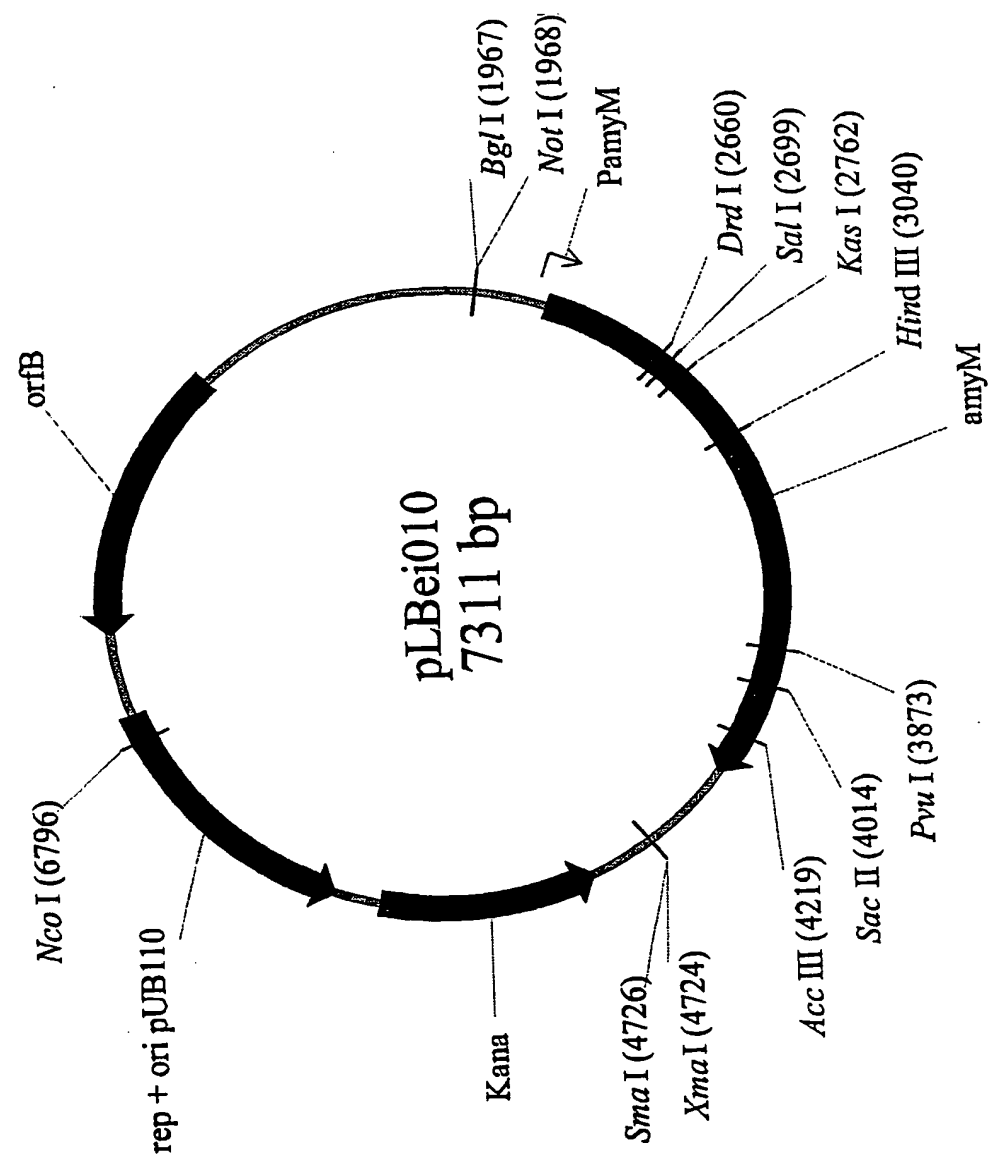
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Fig. 1

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                 20   25   30

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                 35   40   45

1

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        50                  55                  60

ctg ggc gta acg aca atc tgg ttg tcc ccg gtt ttg gac aat ctg gat    336
Leu Gly Val Thr Thr Ile Trp Leu Ser Pro Val Leu Asp Asn Leu Asp
        65                  70                  75

aca ctg gcg ggc acc gat aac acg ggc tat cac gga tac tgg acg cgc    384
Thr Leu Ala Gly Thr Asp Asn Thr Gly Tyr His Gly Tyr Trp Thr Arg
80                  85                  90                  95

gat ttt aaa cag att gag gaa cat ttc ggg aat tgg acc aca ttt gac    432
Asp Phe Lys Gln Ile Glu Glu His Phe Gly Asn Trp Thr Thr Phe Asp
                100                 105                 110

acg ttg gtc aat gat gct cac caa aac gga atc aag gtg att gtc gac    480
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                130                 135                 140

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Ala Glu Gly Gly Ala Leu Tyr Asn Asn Gly Thr Tyr Met Gly Asn Tyr
        145                 150                 155

ttt gat gac gca aca aaa ggg tac ttc cac cat aat ggg gac atc agc    624
Phe Asp Asp Ala Thr Lys Gly Tyr Phe His His Asn Gly Asp Ile Ser
160                 165                 170                 175

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caa tac ctg acc gat gcg gcg gtt caa ttg gta gca cat gga gcg gat    768
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240             245             250             255

gaa tgg tac gga gat gac ccc gga aca gcc aat cat ctg gaa aag gtc    912
Glu Trp Tyr Gly Asp Asp Pro Gly Thr Ala Asn His Leu Glu Lys Val
                260             265             270

cgg tac gcc aac aac agc ggt gtc aat gtg ctg gat ttt gat ctc aac    960
Arg Tyr Ala Asn Asn Ser Gly Val Asn Val Leu Asp Phe Asp Leu Asn
                275             280             285

acg gtg att cga aat gtg ttc ggc aca ttt acg caa acg atg tac gat   1008
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        305             310             315

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320             325             330             335

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Thr Ser Arg Gly Thr Pro Ser Ile Tyr Tyr Gly Thr Glu Gln Tyr Met
                355             360             365

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Ala Gly Gly Asn Asp Pro Tyr Asn Arg Gly Met Met Pro Ala Phe Asp
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agg aac aat gcg gcg atc cag tac ggc acc acc acc cag cgt tgg atc   1344
Arg Asn Asn Ala Ala Ile Gln Tyr Gly Thr Thr Thr Gln Arg Trp Ile
400             405             410             415

aac aat gat gtt tac att tat gaa cgg aaa ttt ttc aac gat gtc gtg   1392
Asn Asn Asp Val Tyr Ile Tyr Glu Arg Lys Phe Phe Asn Asp Val Val
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3

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Phe Thr Leu Ala Pro Gly Ala Val Ser Val Trp Gln Tyr Ser Thr Ser
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gct tca gcg ccg caa atc gga tcg gtt gct cca aat atg ggg att ccg   1632
Ala Ser Ala Pro Gln Ile Gly Ser Val Ala Pro Asn Met Gly Ile Pro
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acc gtg aca ttt ggc gga gtg aca gcg act gtg aaa tcc tgg aca tcc   1728
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Asn Arg Ile Glu Val Tyr Val Pro Asn Met Ala Ala Gly Leu Thr Asp
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560   565   570   575

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cct ccg acc aac ctg ggg gat aag att tac ctg acg ggc aac ata ccg   1920
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4

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5

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Asn Asn Asp Val Tyr Ile Tyr Glu Arg Lys Phe Phe Asn Asp Val Val
    450   455   460

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        580   585   590

Val Lys Val Thr Ala Gly Gly Val Ser Ser Asn Leu Tyr Ser Tyr Asn
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7

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Gln Gly Pro Leu Leu Ala Pro Asn Tyr Pro Asp Trp Phe Tyr Val Phe
            660   665   670

Ser Val Pro Ala Gly Lys Thr Ile Gln Phe Lys Phe Phe Ile Lys Arg
        675   680   685

Ala Asp Gly Thr Ile Gln Trp Glu Asn Gly Ser Asn His Val Ala Thr
    690   695   700

Thr Pro Thr Gly Ala Thr Gly Asn Ile Thr Val Thr Trp Gln Asn
705   710   715

<210> 3
<211> 38
<212> DNA
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<220>
<223> Description of Artificial Sequence: F 188H Primer

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8

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<212> DNA
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<220>
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<210> 7
<211> 36
<212> DNA
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<220>
<223> Description of Artificial Sequence: F 284K Primer

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<210> 8
<211> 35
<212> DNA
<213> Artificial Sequence

<220>
<223> Description of Artificial Sequence: N 327D Primer

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<210> 9
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<212> DNA
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<223> Description of Artificial Sequence: T 288K Primer
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<212> DNA
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<210> 11
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<210> 12
<211> 20
<212> DNA
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<220>
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<210> 13
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<212> DNA
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<210> 14
<211> 22
<212> DNA
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<220>
<223> Description of Artificial Sequence: B 346 Primer

<400> 14
tttgtactcg ttccccgttt gg   22

INTERNATIONAL SEARCH REPORT

International application No.

PCT/DK 99/00088

A. CLASSIFICATION OF SUBJECT MATTER

IPC6: C12N 9/28, A21D 8/04

According to International Patent Classification (IPC) or to both national classification and IPC

B. FIELDS SEARCHED

Minimum documentation searched (classification system followed by classification symbols)

IPC6: C12N

Documentation searched other than minimum documentation to the extent that such documents are included in the fields searched

SE,DK,FI,NO classes as above

Electronic data base consulted during the international search (name of data base and, where practicable, search terms used)

C. DOCUMENTS CONSIDERED TO BE RELEVANT

Category*	Citation of document, with indication, where appropriate, of the relevant passages	Relevant to claim No.
X	WO 9623874 A1 (NOVO NORDISK A/S), 8 August 1996 (08.08.96), see page 30-37, claims 1-2	1-32
	--	
X	WO 9743424 A1 (GENENCOR INTERNATIONAL, INC.), 20 November 1997 (20.11.97), see page 3, lines 21-34, page 12, lines 15-19	1-8
A		7-32
	--	
X	WO 9741213 A1 (NOVO NORDISK A/S), 6 November 1997 (06.11.97), see page 11, lines 5-12	1-8
A		9-32
	--	

[X] Further documents are listed in the continuation of Box C.   [X] See patent family annex.

* Special categories of cited documents:	"T" later document published after the international filing date or priority date and not in conflict with the application but cited to understand the principle or theory underlying the invention
"A" document defining the general state of the art which is not considered to be of particular relevance	
"E" earlier document but published on or after the international filing date	"X" document of particular relevance: the claimed invention cannot be considered novel or cannot be considered to involve an inventive step when the document is taken alone
"L" document which may throw doubts on priority claim(s) or which is cited to establish the publication date of another citation or other special reason (as specified)	
"O" document referring to an oral disclosure, use, exhibition or other means	"Y" document of particular relevance: the claimed invention cannot be considered to involve an inventive step when the document is combined with one or more other such documents, such combination being obvious to a person skilled in the art
"P" document published prior to the international filing date but later than the priority date claimed	"&" document member of the same patent family

Date of the actual completion of the international search	Date of mailing of the international search report
8 June 1999	0 8 -06- 1999
Name and mailing address of the ISA/ Swedish Patent Office Box 5055, S-102 42 STOCKHOLM Facsimile No. + 46 8 666 02 86	Authorized officer Yvonne Siösteen Telephone No. + 46 8 782 25 00

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INTERNATIONAL SEARCH REPORT

C (Continuation). DOCUMENTS CONSIDERED TO BE RELEVANT

Category*	Citation of document, with indication, where appropriate, of the relevant passages	Relevant to claim No.
A	Databas Swissprot, accessionno. P19531, Diderichsen B. et al: "Maltogenicalpha-amylase precursor (EC 3.2.1.133) Glucan 1,4-alpha-maltohydrolase)"; & Fems Microbiol. Lett. 56:53-60 (1988)	1-32
	--	
A	Dialog Information Services, file 5, Biosis Previews, Dialog accession no. 07281496, Biosis accession no. 000090061383, Boel E et al: "Calcium binding in alpha amylases an X-ray diffraction study at 2.1-Aresolution of two enzymes from aspergillus"; & Biochemistry 29 (26). 1990. 6244-6249	1-32
	--	
A	Dialog Information Services, file 5, Biosis Previews, Dialog accession no. 09764348, Biosis accession no. 199598219266, Machius M et al: "Crystal structure of calcium-depleted Bacillus licheniformis alpha-amylase at 2.2 A resolution"; & Journal of Molecular Biology 246 (4):p545-559 1995	1-32
	--	
A	J. Mol. Biol., Volume 229, 1993, Changsoo Chang et al, "Crystallization and Preliminary X-ray Crystallographic Analysis of alpha-Amylase from Bacillus subtilis" page 235 - page 238	1-32
	--	
A	Plant Molecular Biology, Volume 25, 1994, Birte Svensson, "Protein engineering in the alpha-amylase family: catalytic mechanism, substrate specificity, and stability" page 141 - page 157	1-32
	--	

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INTERNATIONAL SEARCH REPORT

Information on patent family members

03/05/99	International application No. PCT/DK 99/00088	

Patent document cited in search report	Publication date	Patent family member(s)	Publication date
WO 9623874 A1	08/08/96	AU 4483496 A BR 9607013 A CA 2211316 A CN 1172501 A EP 0808363 A JP 11500003 T	21/08/96 28/10/97 08/08/96 04/02/98 26/11/97 06/01/99
WO 9743424 A1	20/11/97	AU 2996997 A US 5763385 A	05/12/97 09/06/98
WO 9741213 A1	06/11/97	AU 2692897 A EP 0904360 A	19/11/97 31/03/99